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ANALYTICAL RESULTS REPORT
EMPIRE STATE OIL REFINERY

TDD R8-8509-23

CASE #4242

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SUBMITTED TO: KEITH SCHWAB, FIT-DPO

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DATE SUBMITTED: MAY 6, 1986

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ANALYTICAL RESULTS REPORT
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CASE #4242

I. INTRODUCTION

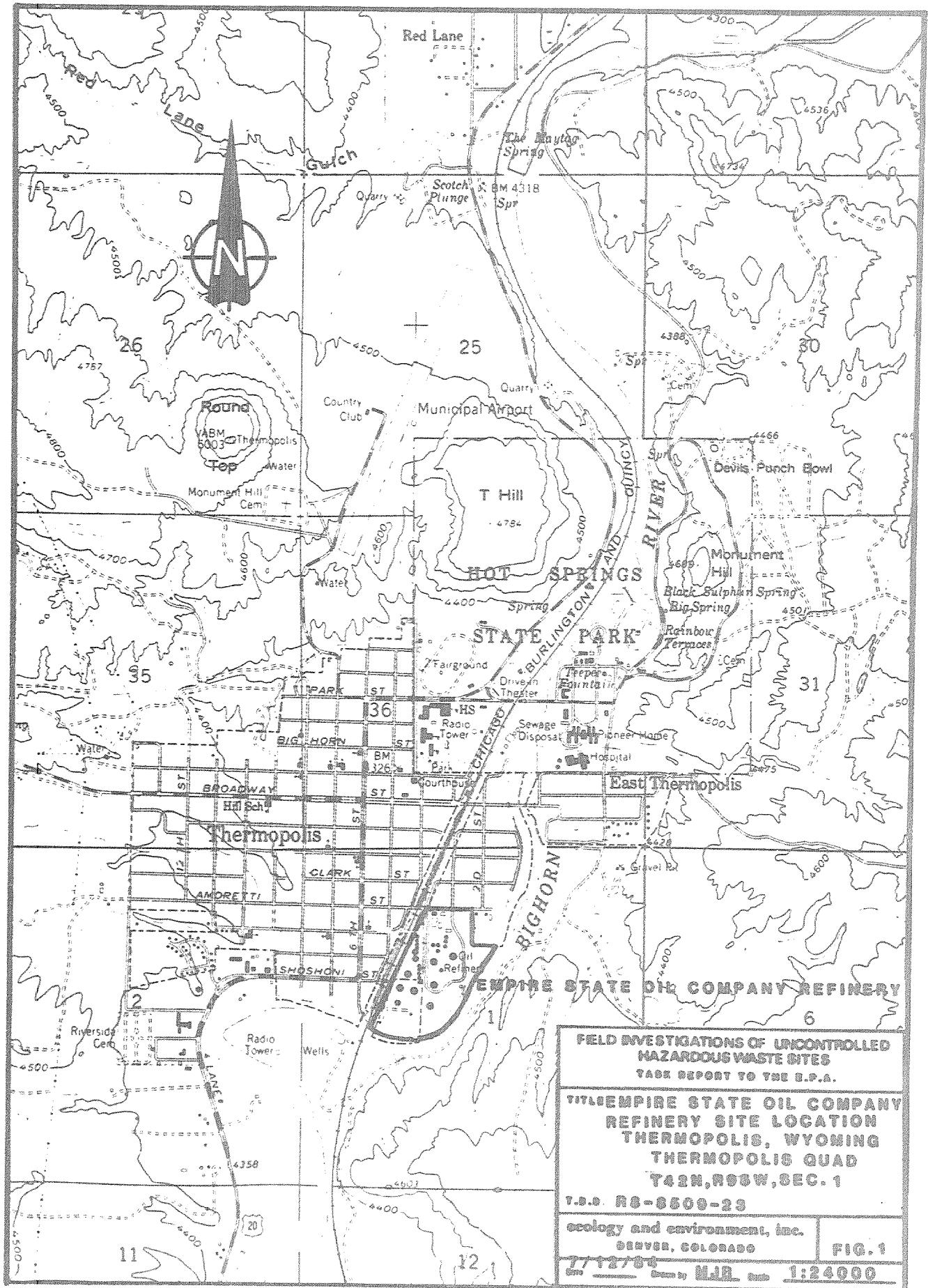
This report was prepared to satisfy the requirements of Technical Directive Document (TDD) R8-8509-23, issued to Ecology and Environment, Inc., Field Investigation Team (E&E FIT) by the Region VIII Environmental Protection Agency (EPA)

The samples discussed in this report were collected by the FIT on April 21, 1985, from Empire State Oil Company, Thermopolis, Wyoming (Figure 1). Two previous reports, the report of sampling activities and the Sampling Plan (R8-8503-20), present discussions regarding project objectives, sample documentation and field observations.

The objective of this investigation was to determine the existence of contamination resulting from activities at the Empire State Oil Company Refinery and the extent of contaminant migration toward the Big Horn River. The FIT also collected data which may be used for the completion of a site inspection (SI).

The Empire State Oil Company Refinery is located at the southeast edge of Thermopolis in central Wyoming. The legal description of the center of the site is NE 1/4, SW 1/4, NW 1/4 of Section 1, T. 42S, R.95W. The site is bounded on the east and south by the Big Horn River, on the west by a railroad spur, and on the north by Amoretti Street.

The refinery was operated by Empire State Oil Company from 1938 until its shutdown in 1969. In August of 1972, Ashland Oil Company acquired all domestic assets of Empire State Oil Company including their refinery installation in Thermopolis, Wyoming. Ashland Oil Company then deeded the property in September of 1972 to the Gottsche



Foundation, which runs a rehabilitation center in Thermopolis. Mr. John Herrin, former president of Empire State Oil Company, stated that the capacity of the refinery was 7,000 barrels per day. Asphalt, fuel oil, diesel fuel, and distillate were produced, but no gasoline was produced because the crude oil that was utilized was too heavy. The facility had an API separator and several pits and ponds that can be seen on a Wyoming Highway Department aerial photograph. A large pond containing water used for fire protection and boiler water remains at the site.

All of the tanks were removed in 1974 by Ashland Oil Company and the site was leveled. Old refinery buildings remain on the north edge of the site, but otherwise the land is covered with vegetation and a few scattered piles of building materials.

II. QUALITY ASSURANCE

Three surface water samples were collected: an upstream, background sample (ES-SW-1), a downstream sample (ES-SW-2), and a sample from the pond (ES-SW-3). Three sediment samples (ES-SE-1, ES-SE-2, and ES-SE-3) were collected from locations corresponding to the surface water samples. Four soil samples were taken: a background sample (ES-SO-1) and three additional samples (ES-SO-2, ES-SO-3 and ES-SO-4) from various on-site locations shown in Figure 2). A field blank water sample (ES-BL-1) was prepared. A field duplicate (ES-SW-30) was taken from the same location as ES-SW-3.

All water, sediment and soil samples collected at Empire State Oil Company were analyzed for Hazardous Substance List (HSL) volatile organics, base/neutral/acid(B/N/A) extractable organics and Task 1 and 2 metals. One sample (ES-SO-4) was designated as medium hazard and was hand-delivered to Rocky Mountain Analytical Laboratory in Denver, Colorado. The remaining samples were shipped to ChemTech for inorganic analyses, and to Environmental Control Technology Corporation for organic analyses.

The organic data package was examined for quality assurance by an E&E FIT data reviewer. The findings are as follows:

1. Holding times were exceeded for samples ES-SW-2 and ES-BL-1 because samples had to be reanalyzed.
2. Several non-SPCC's and CCC's were not within contracts limits for calibration.
3. Methylene chloride, acetone, and 2-butanone were found in the field blank. Qualifications were made on the organic analyses based on these blank observations.
4. Some of the advisory limits on the matrix spike and matrix spike duplicate were exceeded.

The organic data was judged as acceptable for use with the noted qualifications.

The inorganic data package was examined for quality assurance by an E&E FIT reviewer. The data were judged acceptable for use with the qualifications noted below:

1. Many spike recovery requirements were not met. As a result, the data for some elements are flagged with an "r".
2. Low soil results for chromium are flagged with an "*" because of poor duplicate reproducibility.
3. Iron was detected in the blank and some of the samples at similar levels.

The following compounds were analyzed for, but the results are not usable because of poor quality control: vinyl acetate, 2-chloroethylvinylether, benzidine, 2-nitroaniline, benzidine and 3,3'-dichlorobenzidine. In the soil and sediment samples, the data

for antimony, beryllium, chromium, copper, lead, nickel, silver, thallium, and tin were not usable. In the surface water samples, iron mercury, silver, thallium and tin data were not usable.

A number of compounds were detected in the associated laboratory blanks. Where appropriate, these samples were qualified with a "b". A wide variation in the blind duplicate was noted for some organic compounds (acetone and 2-butanone) and metals (aluminum and calcium). The majority of the analyses were in close agreement, however.

III. ANALYTICAL RESULTS

Analytical results for the 1985 sampling effort at Empire State Refinery have been compiled in Tables 1, 2, 3 and 4. Corresponding sample locations are illustrated in Figure 2. Only those organic contaminants detected and only usable data are shown in the tables.

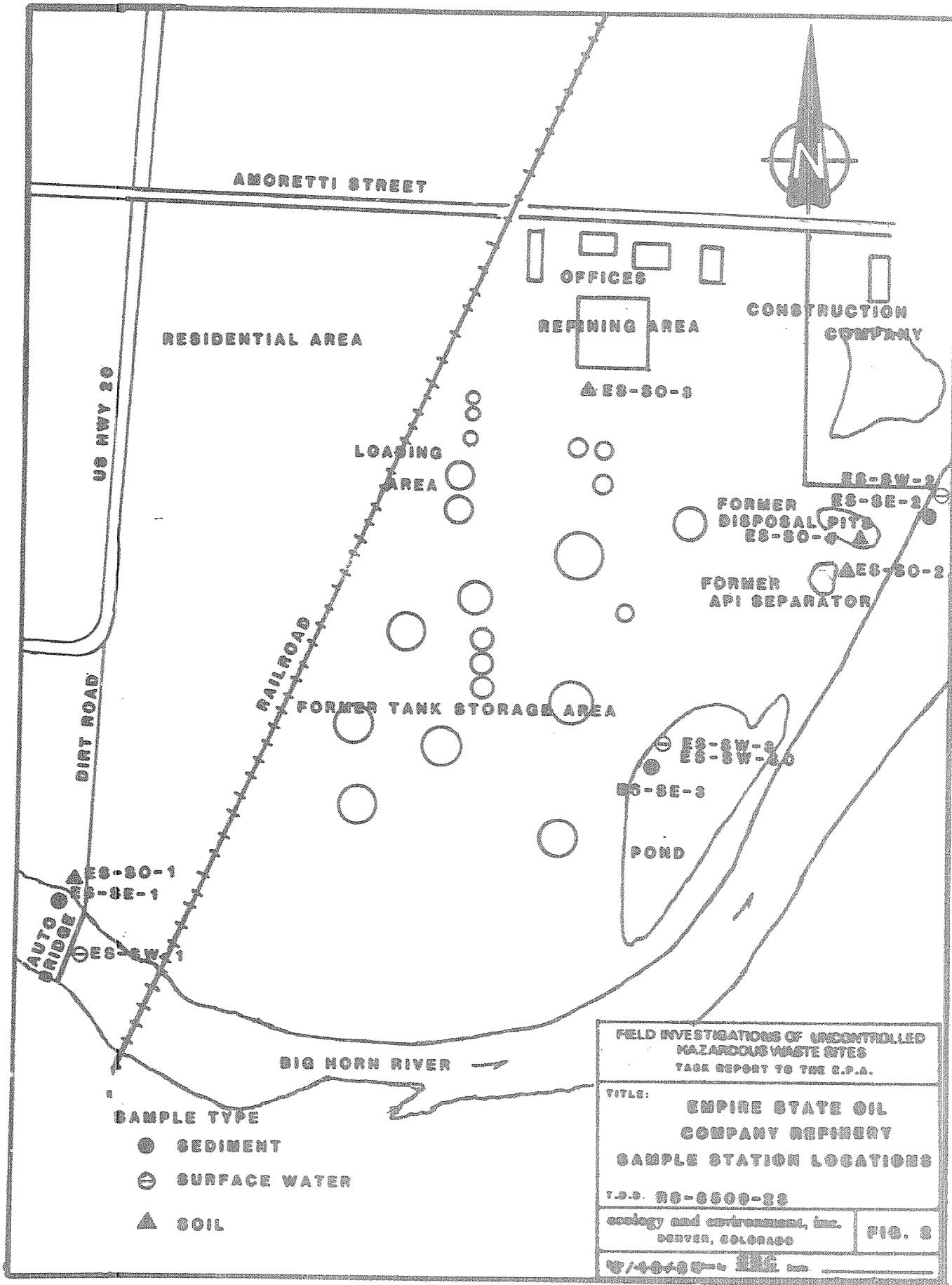
A review of the analytical data allows the following observations and conclusions.

A. SURFACE WATER AND SEDIMENTS

The finding of acetone, methylene chloride and 2-butanone in the surface water and sediment samples is associated with the finding of similar levels in the field decontamination blank and are, therefore, not considered to be associated with site contamination.

Two surface water and two sediment samples were taken from the Big Horn River. A comparison of the downgradient samples, ES-SW-2 and ES-SE-2, with the background samples, ES-SW-1 and ES-SE-1, suggests a release of toluene into the river. The toluene concentration rises from below detection to 57 ug/l in the surface water.

Sediment and surface water samples, ES-SE-3 and ES-SW-3, were collected from the pond shown in Figure 2. A comparison of the pond sediment sample with the background sediment sample (ES-SE-1)



indicates the onsite presence of mercury. The mercury concentration in the pond is 1.7 mg/kg, as compared to below detection in the background sample. There also appears to be a slight release of aluminum and manganese into the pond.

Several tentatively identified compounds, many of which are hydrocarbons, were detected in the sediment samples. A more complete listing is found in Appendix A.

B. SOIL

Soil samples were collected from a background location (ES-SO-1), the former API separator (ES-SO-2), the refinery area (ES-SO-3), and the former disposal pit (ES-SO-4). The sample ES-SO-2 contained mercury at 0.5 mg/kg. Background soil concentrations in the western U.S. have a geometric mean of 0.05 mg/kg and normal range or 0.02 - 0.11 mg/kg (Schacklette, 1984 U.S.G.S. Prof Paper #1270). Toluene was also detected in the sample ES-SO-4 at concentrations far above expected, which may indicate actual presence of high concentrations in the soil.

IV. DISCUSSION

Thermopolis and the surrounding communities (population approximately 5,000) utilize surface water from the Big Horn River at the southwestern end of the Refinery. This source is augmented by ground water from a well field approximately 1,600 feet west-southwest of the Refinery during the summer equal to 50% of total usage.

The source of the municipal ground water supply is the Quaternary alluvial aquifer. As described by Berry and Littleton (1961) and Libra et al (1981), the hydrologic characteristics of this aquifer are:

Specific Capacity	0.3 - 70 gpm/ft
Permeability	2,200 - 4,400 gpd/ft ²
Transmissivity	200 - 80,000 gpd/ft

Recharge occurs primarily from direct penetration of precipitation,

and seepage from irrigation canals, surface streams and depressions. Discharge occurs mainly by evapotranspiration, stream gains and discharge from wells seeps and springs. Recharge of the aquifer occurs primarily from the river during the spring months and the process is reversed during the fall.

No data were available on aquifer storage coefficients and well drawdowns. The piezometric surfaces were at approximately 8 feet and the average depth of wells at 30 feet.

V. CONCLUSIONS

The above analytical results indicate a possible release of toluene into the Big Horn River from Empire State Oil Refinery. Mercury concentrations from the API separator soil sample were also elevated. However, the mercury results for water were of such poor quality that no conclusion could be reached from those data.

The FIT recommends a resampling of this site for the following reasons:

- a. More data are needed for a better assessment of ground water contaminant release.
- b. Approximately 1,600' to the south of ES-SO-4 is an active well field used to supply the community of Thermopolis and surrounding area (approximately 5,000 residents).
- c. More completely assess the extent of the surface water release to the Big Horn River.

REFERENCES

Berry, D. W. and Littleton, R.T., Geology and Ground Water Resources of the Owl Creek Area, Hot Springs County, Wyoming, U.S.G.S. Water Supply Paper 1519, 1961.

Libra, R. et al, Occurrence and Characteristics of Ground Water in the Bighorn Basin, Wyoming. Water Resources Research Institute, University of Wyoming, June 1981.

TABLE 1

ORGANIC ANALYTICAL RESULTS, WATER SAMPLES (ug/l)
 EMPIRE STATE OIL REFINERY, THERMOPOLIS, WYOMING
 CASE #4242

	H1838 ES-SW-1	H1839 ES-SW-2	H1840 ES-SW-3	H1841 ES-SW-30	H1842 ES-BL-1
Methylene Chloride	1600j	300j	12j	6j	22j
Acetone	82j	38j	140j		1400j
2-Butanone	14j	420j	200j		22j
Bis(2-ethylhexyl)phthalate	7j	2j	5j	5j	4j
Toluene		57	7		
Di-N-Butylphthalate			3		4

j - The associated numerical value is an estimated quantity because the amount detected is below the required limits or because quality control criteria were not met.

TABLE 2

ORGANIC ANALYTICAL RESULTS, SOILS ND SEDIMENTS (ug/kg)
 EMPIRE STATE OIL REFINERY, THERMOPOLIS, WYOMING
 CASE #4242

	H1843 ES-SE-1	H1845 ES-SE-2	H1849 ES-SE-3	H1844 ES-SO-1	H1846 ES-SO-2	H1847 ES-SO-3	
Methylene Chloride	15j	40j	72j	15j	15j	10j	2500b
Acetone	22j	33j	37j	3j		3j	29000b
2-Butanone	6j	5j					
Di-N-Butylphthalate	700ub	660ub		540ub		450ub	
Bis (2-ethylhexyl) phthalate	1,100j	840j		290j		380j	
Toluene			3jb		37b	7ub	1500b

b - Compound was detected in the blank. Quantity reported is 5 X the amount found in the blank (10 for methylene chloride, acetone, toluene, nd phthalates).

ub - Estimated sample quantitation limit increased. Amount found in sample reported. Compound detected at 5 X the amount in blank (10 X for methylene chloride, acetone, toluene and phthalates).

j - The associated numerical value is an estimated quantity because the amount detected is below the required limits or because quality control criteria were not met.

jb - The value is an estimated amount detected below required limits and also detected in the blank.

TABLE 3

INORGANIC ANALYTICAL RESULTS, WATER SAMPLES (ug/l)
 EMPIRE STATE OIL REFINERY, THERMOPOLIS, WYOMING
 CASE #4242

	MBH243 ES-SW-1	MBH249 ES-SW-2	MBH244 ES-SW-3	MBH245 ES-SW-30	MBH246 ES-BL-1
Aluminum	[147]	50u	1763	50u	50u
Antimony	50u	50u	50u	50u	50u
Arsenic	10u	10u	10u	10u	10u
Barium	50u	50u	50u	50u	50u
Beryllium	5u	5u	5u	5u	5u
Cadmium	5u	5u	5u	5u	5u
Calcium	61,580	50,890	71,040	46,290	[878]
Chromium	10u	10u	10u	10u	10u
Cobalt	20u	20u	20u	20u	20u
Copper	[23]	20u	[21]	20u	20u
Iron	[82]	20u	2,519	[83]p	[86]r
Lead	5u	5u	5u	5u	5u
Magnesium	20,150	17,260	87,900	86,870	150u
Manganese	15u	15u	247	147	15u
Mercury	0.2ur	0.2u	0.2u	0.2ur	0.2u
Nickel	[21]	20u	20u	20u	20u
Potassium	[2180]	[1762]	9297	8590	1000u
Selenium	5u	5u	5u	5u	5u
Silver	10u	10ur	10ur	10ur	10ur
Sodium	60,230	50,090	297,300	299,200	[1997]
Thallium	10u	10ur	10utr	10urfr	10ur
Tin	30u	30ur	300ur	300ur	30ur
Vanadium	20u	20u	20u	20u	20u
Zinc	32	[19]	28	41	20u
Cyanide	N/R	N/R	N/R	N/R	N/R

u - Material was analyzed for, but not detected. Value is the estimated sample quantitation limit.

[]- Value is below required reporting value.

R - Spike recovery did not meet required recovery limits.

TABLE 4
 INORGANIC ANALYTICAL RESULTS, SOILS AND SEDIMENTS (mg/kg)
 EMPIRE STATE OIL REFINERY, THERMOPOLIS, WYOMING
 CASE #4242

	MHB247 ES-SE-1	MHB248 ES-SE-2	MHB253 ES-SE-3	MHB242 ES-SO-1	MHB250 ES-SO-2	MHB251 ES-SO-3
Aluminum	11226	7385	6192	18646	8255	12031
Antimony	40ur	32ur	47ur	29ur	32ur	25ur
Arsenic	8.0us	6.4u	9.4u	58	9.5	9.6
Barium	201	165	666	340	219	252
Beryllium	4.0ur	3.2ur	4.7ur	2.9ur	3.2ur	2.5ur
Cadmium	4.0u	4.3	4.7u	2.9u	3.2u	2.5u
Calcium	57580	47270	248214	52970	29594	40861
Chromium	29r*	14r*	14r*	39r*	26r*	31r*
Cobalt	16u	13u	19u	12u	13u	10u
Copper	16ur	13ur	19u	32r	13ur	14r
Iron	14503	10371	10226	24131	12957	15912
Lead	17r	5.9sr	128r	29r	3.2ur	16r
Magnesium	15970	11714	9492	17625	8141	13045
Manganese	370	353	1284	596	308	422
Mercury	0.16u	0.13u	1.7	0.12u	0.51	0.10u
Nickel	[19]r	13u	19ur	31r	[16]r	24r
Potassium	[2314]	[1395]	[1252]	5130	[1409]	[2048]
Selenium	4.0u	3.2u	4.7u	2.9u	3.2u	2.5u
Silver	8.6	6.4ur	9.4ur	5.8ur	6.3ur	5.1ur
Sodium	[941]	[1189]	[1902]	[996]	761u	[976]
Thallium	8.0ur	6.4ur	9.4ur	5.8ur	6.3ur	5.1ur
Tin	24ur	19ur	28ur	18ur	19ur	15ur
Vanadium	[29]	[22]	[24]	43	[29]	33
Zinc	37	18	23	119	19	32

u - Material was analyzed for, but not detected. Value is the estimated sample quantitation limit.

[] - Value is below required reporting limit

r - Spike recovery did not meet required recovery limits.

* - Duplicate results exceed R.P.D. limit.

APPENDIX A
QUALITY ASSURANCE REPORTS

REGION VIII SUMMARY OF DATA QUALITY ASSURANCE REVIEW

Case No. 4242 Project No. _____
Site Empire Refinery
Contractor Laboratory Chemtech
Data Reviewer dm Roberts Date of Review 7/8/95
Sample Matrix 5 water + 6 soil Metal Analysis

Sample No. MHB 242 MHB 247 MHB 253

243 248 _____
244 249 _____
245 250 _____
246 251 _____

- () Data are acceptable for use
(✓) Data are acceptable for use with qualification noted above below
() Data are preliminary - pending action or verification
() Data are unacceptable

Action required by DPO?

No ✓ Yes _____ Following items require action _____

Action required by Project Officer (PO)?

No ✓ Yes _____

Form A

Inorganic Data Completeness Checklist

- Inorganic analysis data sheets
- Instrument Detection limits
- Duplicate results
- Spike results
- ICP interference check sample
- Blank results
- Raw data for calibration standards
- Raw data for blanks
- Raw data for samples
- Raw data for duplicates
- Raw data for spikes
- Initial calibration and calibration verification results
- Continuing calibration verification
- Traffic Reports

Form B

All inorganic standards were within specified contract limits. - - - - -

Yes

No

Comments:

All inorganic detection limits met the contract requirements.

Yes

No

Comments:

All matrix spike requirements were met.

Yes

No

Comments:

Raw Water:

Iron	14%
Manganese	20%
Silver	0%
Ruthenium	50%
Tin	0%

Raw Soil:	Antimony	57%
	Beryllium	126%
	Chromium	127%
	Copper	128%
	Lead	264%
	Nickel	127%
	Silicon	24%
	Thallium	148%
		200%

The interference check sample was run twice per eight hour shift. No massive interferences were present.

Yes

No

Comments:

Form C

A blank was run with every twenty samples or less per case.

Yes No

How many elements were detected above the required detection limit? 0

How many elements were detected at greater than one half the amount detected in any sample?: 1

Comments: Iron detected in calibration and preparation blank at ~60 µg/l. Iron was detected in some low water samples at ~80 µg/l.

A duplicate sample was run with every twenty or fewer samples of a similar matrix, or one per case, whichever is more frequent.

Yes No

The RPD's were tabulated.

Yes No

Comments:

Low Soil: Chromium 48% RPD

All holding times were met.

Yes No

Comments:

Form D

Initial calibration data were reviewed. Initial calibration data were included in the package and met all contract requirements.

Yes

No

Comments:

Continuing calibration data were reviewed and this data met all contract requirements.

Yes

No

Comments:

Following are our findings:

Many spike recovery requirements were not met, as a result the counts for some elements are flagged with an "R". Low soil results for chromium are flagged with an "+" because of poor duplicate reproducibility. It is noted on form C that iron was detected on the blanks and some of the samples at similar levels.

Form I

U.S. EPA Contract Laboratory Program
Sample Management Office
P.O. Box 818 - Alexandria, VA 22313
703/557-2490 FTS: 8-557-2490

EPA Sample No.
MHB 245

Date 5-30-85

INORGANIC ANALYSIS DATA SHEET

LAB NAME CHEMTECH CONSULTING GROUP

CASE NO. 4242

SOW NO. 784

LAB SAMPLE ID. NO. 62-473-01

QC REPORT NO. 473

Elements Identified and Measured

Concentrations:

Low /

Medium /

Matrix: Water /

Soil /

Sludge /

Other /

(ug/L or mg/kg dry weight (Circle One))

1. Aluminum	<u>147</u> P.
2. Antimony	<u>50</u> U.P.
3. Arsenic	<u>10</u> U.P. UF.
4. Barium	<u>50</u> U.P.
5. Beryllium	<u>5</u> U.P.
6. Cadmium	<u>5</u> U.P.
7. Calcium	<u>61580</u> P.
8. Chromium	<u>10</u> U.P.
9. Cobalt	<u>20</u> U.P.
10. Copper	<u>23</u> P.
11. Iron	<u>182</u> P. R
12. Lead	<u>5</u> U.F.
Glycide	<u>NR</u>

13. Magnesium	<u>20150</u> P.
14. Manganese	<u>15</u> U.P.
15. Mercury	<u>0.2</u> U.R
16. Nickel	<u>217</u> P.
17. Potassium	<u>2180</u> P.
18. Selenium	<u>5</u> U.F.
19. Silver	<u>10</u> U.P. R
20. Sodium	<u>60230</u> P.
21. Thallium	<u>10</u> U.F. R
22. Tin	<u>30</u> U.F. R
23. Vanadium	<u>20</u> U.P.
24. Zinc	<u>32</u> P.

Percent Solids (%)

Footnotes: For reporting results to EPA, standard result qualifiers are used as defined on Cover Page. Additional flags or footnotes explaining results are encouraged. Definition of such flags must be explicit and contained on Cover Page, however.

Comments:

Lab Manager L. H. Keddy

Form I

U.S. EPA Contract Laboratory Program
Sample Management Office
P.O. Box 810 - Alexandria, VA 22313
703/537-2490 PTS: 8-537-2490

EPA Sample No.

MHB 244

Date 5-30-PS

INORGANIC ANALYSIS DATA SHEET

LAB NAME CHEMTECH CONSULTING GROUP

CASE NO. 4242

SOW NO. 784

LAB SAMPLE ID. NO. G2-473-02

QC REPORT NO. 473

Elements Identified and Measured

Concentration:

Low /

Medium /

Matrix: Water /

Soil /

Sludge /

Other /

ug/L or ug/kg dry weight (Circle One)

1. Aluminum 1763 P.
2. Antimony 50 UP.
3. Arsenic 10 UP. UF.
4. Barium 50 UP.
5. Beryllium 5 UP.
6. Cadmium 5 UP.
7. Calcium 71040 P.
8. Chromium 10 UP.
9. Cobalt 20 UP.
10. Copper [21] P.
11. Iron 2519 P.R
12. Lead 5 UF.

Cyanide NR

13. Magnesium 87900 P.
14. Manganese 247 P.
15. Mercury 0.2 U R
16. Nickel 20 UP.
17. Potassium 9297 P.
18. Selenium 5 UF.
19. Silver 10 UP. R
20. Sodium 297300 P.
21. Thallium 10 UF. R
22. Tin 300 UF. R
23. Titanium 20 UP.
24. Zinc 28 P.

Percent Solids (%)

Footnotes: For reporting results to EPA, standard result qualifiers are used as defined on Cover Page. Additional flags or footnotes explaining results are encouraged. Definitions of such flags must be explicit and contained on Cover Page, however.

Comments:

Lab Manager



Form I

U.S. EPA Contract Laboratory Program
Sample Management Office
P.O. Box 818 - Alexandria, VA 22313
703/537-2490 FTS: 8-537-2490

EPA Sample No.

MHB 245

Date 5-20 85

INORGANIC ANALYSIS DATA SHEET

LAB NAME CHEMTECH CONSULTING GROUP

CASE NO. 4242

SOW NO. 784

LAB SAMPLE ID. NO. G2-473-03

QC REPORT NO. 473

Elements Identified and Measured

Concentration:

Lov ✓

Medium _____

Matrix: Water /

Soil _____

Sludge _____

Other _____

(ug/L or ug/kg dry weight (Circle One))

1. Aluminum 50 UP.
2. Antimony 50 UP.
3. Arsenic 10 UP. UF
4. Barium 50 UP.
5. Beryllium 5 UP.
6. Cadmium 5 UP.
7. Calcium 46290 P.
8. Chromium 10 UP.
9. Cobalt 20 UP.
10. Copper 20 UP.
11. Iron 1837 P. R
12. Lead 5 UF.
Cyanide NR

13. Magnesium 86870 P.
14. Manganese 147 P.
15. Mercury 0.2 UP R
16. Nickel 20 UP.
17. Potassium 8590 P.
18. Selenium 5 UF.
19. Silver 10 UP. R
20. Sodium 299200 P.
21. Thallium 10 UF. R
22. Tin 300 UF. R
23. Titanium 20 UP.
24. Zinc 41 P.

Percent Solids (%)

Footnotes: For reporting results to EPA, standard result qualifiers are used as defined on Cover Page. Additional flags or footnotes explaining results are encouraged. Definition of such flags must be explicit and contained on Cover Page, however.

Comments:

Lab Manager

E.J.P.
(Hedvat)

Form I

U.S. EPA Contract Laboratory Program
 Sample Management Office
 P.O. Box 810 - Alexandria, VA 22313
 703/557-2490 FTS: 8-557-2490

EPA Sample No.

MHB 246

Date 5-30-85

INORGANIC ANALYSIS DATA SHEET

LAB NAME CHEMTECH CONSULTING GROUP

CASE NO. 4242

SOW NO. 784

LAB SAMPLE ID. NO. G2-473-04

QC REPORT NO. 473

Elements Identified and Measured

Concentration:

Low ✓

Medium _____

Matrix: Water /

Soil _____

Sludge _____

Other _____

(ug/L or ug/kg dry weight (Circle One))

1. <u>Aluminum</u>	<u>50 UP.</u>
2. <u>Antimony</u>	<u>50 UP.</u>
3. <u>Arsenic</u>	<u>10 UP. UF</u>
4. <u>Barium</u>	<u>50 UP.</u>
5. <u>Beryllium</u>	<u>5 UP.</u>
6. <u>Cadmium</u>	<u>5 UP.</u>
7. <u>Calcium</u>	<u>[878] P.</u>
8. <u>Chromium</u>	<u>10 UP.</u>
9. <u>Cobalt</u>	<u>20 UP.</u>
10. <u>Copper</u>	<u>20 UP.</u>
11. <u>Iron</u>	<u>[86] P. R</u>
12. <u>Lead</u>	<u>5 UF.</u>
<u>Cyanide</u>	<u>NR</u>

13. <u>Magnesium</u>	<u>150 UP.</u>
14. <u>Manganese</u>	<u>15 UP.</u>
15. <u>Mercury</u>	<u>0.2 UP.</u>
16. <u>Nickel</u>	<u>20 UP.</u>
17. <u>Potassium</u>	<u>1000 UP.</u>
18. <u>Selenium</u>	<u>5 UF.</u>
19. <u>Silver</u>	<u>10 UP. R</u>
20. <u>Sodium</u>	<u>[1997] P.</u>
21. <u>Thallium</u>	<u>10 UF. R</u>
22. <u>Tin</u>	<u>30 UF. R</u>
23. <u>Vanadium</u>	<u>20 UP.</u>
24. <u>Zinc</u>	<u>20 UP.</u>

Percent Solids (%)

Footnotes: For reporting results to EPA, standard result qualifiers are used as defined on Cover Page. Additional flags or footnotes explaining results are encouraged. Definition of such flags must be explicit and contained on Cover Page, however.

Comments:

Lab Manager

L. Hedrick

Form I

U.S. EPA Contract Laboratory Program
Sample Management Office
P.O. Box 816 - Alexandria, VA 22313
703/537-2490 FTS: 8-537-2490

EPA Sample No.

MMB 249

Date 5-30-85

INORGANIC ANALYSIS DATA SHEET

LAB NAME CHEMTECH CONSULTING GROUP

CASE NO. 4242

SOW NO. 784

LAB SAMPLE ID. NO. G2-473-05

QC REPORT NO. 473

Elements Identified and Measured

Concentration:

Low /

Medium /

Matrix: Water /

Soil /

Sludge /

Other /

ug/L or ug/kg dry weight (Circle One)

1. Aluminum 50 UP.
2. Antimony 50 UP.
3. Arsenic 10 UP. UF.
4. Barium 50 UP.
5. Beryllium 5 UP.
6. Cadmium 5 UP.
7. Calcium 50890. P.
8. Chromium 10 UP.
9. Cobalt 20 UP.
10. Copper 20 UP.
11. Iron 20 UP. R
12. Lead 5 UF.

Cyanide NR

13. Magnesium 17260 P.
14. Manganese 15 UP.
15. Mercury 0.2 UP R
16. Nickel 20 UP.
17. Potassium [1762] P.
18. Selenium 5 UF.
19. Silver 10 UP. R
20. Sodium 50090 P.
21. Thallium 10 UF. R
22. Tin 30 UP. R
23. Titanium 20 UP.
24. Zinc 1197 P.

Percent Solids (2)

Footnotes: For reporting results to EPA, standard result qualifiers are used as defined on Cover Page. Additional flags or footnotes explaining results are encouraged. Definitions of such flags must be explicit and contained on Cover Page, however.

Comments:

Lab Manager

L. Hejna

Form I

U.S. EPA Contract Laboratory Program
Sample Management Office
P.O. Box 818 - Alexandria, VA 22313
703/537-2490 PTS: 8-537-2490

EPA Sample No.

MHS 242

Date 5-30-85

INORGANIC ANALYSIS DATA SHEET

LAB NAME CHEMTECH CONSULTING GROUP

CASE NO. 4242

SOW NO. 784

QC REPORT NO. 473

LAB SAMPLE ID. NO. G2-473-06

Elements Identified and Measured			
Concentrations:	Low /	Medium	
Matrix: Water	Soil /	Sludge	Other

ug/L or ug/kg dry weight (Circle One)

1. <u>Aluminum</u>	<u>18646 P</u>	13. <u>Magnesium</u>	<u>17625 P</u>
2. <u>Antimony</u>	<u>29 UPR</u>	14. <u>Manganese</u>	<u>596 P</u>
3. <u>Arsenic</u>	<u>58 DP UF</u>	15. <u>Mercury</u>	<u>0.12 U</u>
4. <u>Barium</u>	<u>340 P</u>	16. <u>Nickel</u>	<u>31 P R</u>
5. <u>Beryllium</u>	<u>2.9 UPR</u>	17. <u>Potassium</u>	<u>5130 P</u>
6. <u>Cadmium</u>	<u>2.9 UP</u>	18. <u>Selenium</u>	<u>2.9 UF</u>
7. <u>Calcium</u>	<u>52970 P</u>	19. <u>Silver</u>	<u>5.8 UPR</u>
8. <u>Chromium</u>	<u>39 P R*</u>	20. <u>Sodium</u>	<u>[996] P</u>
9. <u>Cobalt</u>	<u>12 UP</u>	21. <u>Thallium</u>	<u>5.8 UF R</u>
10. <u>Copper</u>	<u>32 P R</u>	22. <u>Tin</u>	<u>18 UF R</u>
11. <u>Iron</u>	<u>24131 P</u>	23. <u>Titanium</u>	<u>43 P</u>
12. <u>Lead</u>	<u>29 F. R</u>	24. <u>Zinc</u>	<u>119 P</u>
Cyanide		Percent Solids (%)	<u>85.7</u>

Footnotes: For reporting results to EPA, standard result qualifiers are used as defined on Cover Page. Additional flags or footnotes explaining results are encouraged. Definition of such flags must be explicit and contained on Cover Page, however.

Comments:

Lab Manager

E. Heidrich

Form I

U.S. EPA Contract Laboratory Program
Sample Management Office
P.O. Box 818 - Alexandria, VA 22313
703/597-2490 PTS: 8-597-2490

EPA Sample No.

MNB 247

Date 5-20-85

INORGANIC ANALYSIS DATA SHEET

LAB NAME CHEMTECH CONSULTING GROUP

CASE NO. 4242

SOW NO. 784

LAB SAMPLE ID. NO. G2-473-07

QC REPORT NO. 473

Elements Identified and Measured

Concentration:

Low /

Medium /

Matrix: Water

Soil /

Sludge /

Other /

ug/L or (ug/kg dry weight) (Circle One)

1. Aluminum 11226 F
2. Antimony 40 UPR
3. Arsenic 8.0 UFS
4. Barium 201 P
5. Beryllium 4.0 UPR
6. Cadmium 4.0 UP
7. Calcium 575.80 P
8. Chromium 29 P R*
9. Cobalt 16 UP
10. Copper 16 UPR
11. Iron 14503 P
12. Lead 17 F. R

Cyanide /

13. Magnesium 15970 P
14. Manganese 370 P
15. Mercury 0.16 U
16. Nickel [19] P R
17. Potassium [2314] P
18. Selenium 4.0 UF
19. Silver 8.6 P R
20. Sodium [941] P
21. Thallium 8.0 UF R
22. Tin 24 UF R
23. Vanadium [29] P
24. Zinc 37 P

Percent Solids (%) 62.4

Footnotes: For reporting results to EPA, standard result qualifiers are used as defined on Cover Page. Additional flags or footnotes explaining results are encouraged. Definition of such flags must be explicit and contained on Cover Page, however.

Comments:

Lab Manager

E Hedrick

Form I

Environmental Monitoring Program
U.S. Postal Service
Post Office Box 12000, Falls Church, VA 22313
Telephone 703-537-2490

SEPA
PCB

Date

ENVIRONMENTAL DATA SHEET

WATER GROUP

CASE NO. 0

SOIL NO.

784

10-27-74/13-08

REPORT NO.

Constituents Analyzed and Measured

Condition:

Low ✓

Medium

Soil Type:

Soil ✓

Sludge

Water _____

Metals (Circle One)

1. Arsenic
2. Barium
3. Cadmium
4. Chromium
5. Cobalt
6. Copper
7. Lead
8. Manganese
9. Mercury
10. Nickel
11. Palladium
12. Platinum
13. Silver
14. Sodium
15. Tellurium
16. Tin
17. Uranium
18. Zinc

Other Solids (2)

Indicate other substances being analyzed because of their presence in the sample or for analytical convenience of such items as glass, plastic, or rubber.

Sample

Storage

Form I

U.S. EPA Contract Laboratory Program
Sample Management Office
P.O. Box 818 - Alexandria, VA 22313
703/557-2490 FTS: 8-557-2490

EPA Sample No.

MHB 250

Date 5-30-95

INORGANIC ANALYSIS DATA SHEET

LAB NAME CHEMTECH CONSULTING GROUP

CASE NO. 4242

SOW NO. 784

LAB SAMPLE ID. NO. G2-473-09

QC REPORT NO. 473

Elements Identified and Measured

Concentration:

Low /

Medium /

Matrix: Water

Soil /

Sludge /

Other /

ug/L or mg/kg dry weight (Circle One)

1. <u>Aluminum</u>	<u>8255 P</u>	13. <u>Magnesium</u>	<u>8141 P</u>
2. <u>Antimony</u>	<u>32 UPR</u>	14. <u>Manganese</u>	<u>308 P</u>
3. <u>Arsenic</u>	<u>9.5 F</u>	15. <u>Mercury</u>	<u>0.51</u>
4. <u>Barium</u>	<u>219 P</u>	16. <u>Nickel</u>	<u>[167] P R</u>
5. <u>Beryllium</u>	<u>3.2 UPR</u>	17. <u>Potassium</u>	<u>[1409] P</u>
6. <u>Cadmium</u>	<u>3.2 UP</u>	18. <u>Selenium</u>	<u>3.2 UF</u>
7. <u>Calcium</u>	<u>29594 P</u>	19. <u>Silver</u>	<u>6.3 UPR</u>
8. <u>Chromium</u>	<u>26 P R</u>	20. <u>Sodium</u>	<u>761 UP</u>
9. <u>Cobalt</u>	<u>13 UP</u>	21. <u>Thallium</u>	<u>6.3 UFR</u>
10. <u>Copper</u>	<u>13 UPR</u>	22. <u>Tin</u>	<u>19 UFR</u>
11. <u>Iron</u>	<u>12957 P</u>	23. <u>Vanadium</u>	<u>[297] P</u>
12. <u>Lead</u>	<u>3.2 UFS R</u>	24. <u>Zinc</u>	<u>19 P</u>
Cyanide		Percent Solids (%)	<u>78.8</u>

Footnotes: For reporting results to EPA, standard result qualifiers are used as defined on Cover Page. Additional flags or footnotes explaining results are encouraged. Definition of such flags must be explicit and contained on Cover Page, however.

Comments:

Lab Manager



Form I

U.S. EPA Contract Laboratory Program
Sample Management Office
P.O. Box 818 - Alexandria, VA 22313
703/537-2490 FTS: 8-537-2490

EPA Sample No.

MHB 251

Date 5-30-85

INORGANIC ANALYSIS DATA SHEET

LAB NAME CHEMTECH CONSULTING GROUP

CASE NO. 4342

SOW NO. 784

LAB SAMPLE ID. NO. G2-473-10

QC REPORT NO. 473

Elements Identified and Measured

Concentration:

Low /

Medium /

Matrix: Water

Soil /

Sludge /

Other /

ug/L or mg/kg dry weight (Circle One)

1. Aluminum 12031 P
2. Antimony 25 UPR
3. Arsenic 9.6 F
4. Barium 352 P
5. Beryllium 2.5 UPR
6. Cadmium 2.5 UP
7. Calcium 40861 P
8. Chromium 31 P R*
9. Cobalt 10 UP
10. Copper 14 P R
11. Iron 15912 P
12. Lead 16 F. R
Gypsum

13. Magnesium 13045 P
14. Manganese 422 P
15. Mercury 0.10 U
16. Nickel 24 P R
17. Potassium [2045] P
18. Selenium 2.5 UF
19. Silver 5.1 UPR
20. Sodium [976] P
21. Thallium 5.1 UF R
22. Tin 15 UF R
23. Titanium 33 P
24. Zinc 32 P
Percent Solids (%) 98.7

Footnotes: For reporting results to EPA, standard result qualifiers are used as defined on Cover Page. Additional flags or footnotes explaining results are encouraged. Definition of such flags must be explicit and contained on Cover Page, however.

Comments:

Lab Manager

L. Hart

Form I

U.S. EPA Contract Laboratory Program
Sample Management Office
P.O. Box 810 - Alexandria, VA 22313
703/537-2490 FTS: 8-537-2490

EPA Sample No.

MHB 253

Date 5-30-85

INORGANIC ANALYSIS DATA SHEET

LAB NAME CHEMTECH CONSULTING GROUP

CASE NO. 4242

SOW NO. 784

LAB SAMPLE ID. NO. G2-473-11

QC REPORT NO. 473

Elements Identified and Measured

Concentration:

Low /

Medium /

Matrix: Water

Soil /

Sludge /

Other /

ug/L or ng/kg dry weight (Circle One)

1. Aluminum 619.2 P
2. Antimony 47 UPR
3. Arsenic 9.4 UF.
4. Barium 1166 P
5. Beryllium 4.7 UPR
6. Cadmium 4.7 UP
7. Calcium 248.34 P
8. Chromium 14 P R *
9. Cobalt 19 UP
10. Copper 19 UPR
11. Iron 10.226 P
12. Lead 128 F. R

Cyanide /

13. Magnesium 9492 P
14. Manganese 1284 P
15. Mercury 1.7
16. Nickel 19 UPR
17. Potassium [1252] P
18. Selenium 4.7 UF.
19. Silver 9.4 UPR
20. Sodium [1902] P
21. Thallium 9.4 UF. R
22. Tite 28 UF. R.
23. Vanadium [24] P
24. Zinc 23 P

Percent Solids (%) 53.2

Footnotes: For reporting results to EPA, standard result qualifiers are used as defined on Cover Page. Additional flags or footnotes explaining results are encouraged. Definition of such flags must be explicit and contained on Cover Page, however.

Comments: _____

Lab Manager

E. Hecht

SECTION VIII SUMMARY OF DATA QUALITY ASSURANCE REVIEW

Case No. No Case Number* Project No. R8-8602-11
Site Empire Oil Refinery
Contractor Laboratory Rocky Mountain Analytical
Data Reviewer Randy Greaves Date of Review 2/25/85
Sample Matrix 1 Soil - oily sample

Sample No. ES-50-4 Refining Area.

- Data are acceptable for use
- Data are acceptable for use with qualification noted above
- Data are preliminary - pending action or verification
- Data are unacceptable

Action required by DPO?

No Yes Following items require action

Action required by Project Officer (PO)?

No Yes

* This was not a regular analytical service, and did not involve a laboratory request by SMO.

Following are our findings:

One sample from Empire Oil Refinery and two samples from Ellerby's Oil Refinery were analyzed together. The QA/QC was the same for all three of these samples and, therefore, the data quality assurance review forms will be identical for both the Ellerby's & Empire review. This was an SAS laboratory request for Oily waste. VOA & BNA compounds were analyzed for, but pesticides / PCB's were not analyzed. The laboratory did perform all of the specified requirements of the SAS request and the overall quality of the data produced was good. The % difference between the initial and continuing calibration was out of control for several of the HSL compounds but since none of these compounds were observed in the samples, the effect is probably negligible. The Oily matrix was effectively removed from the samples by the extraction procedure and the quality control was such that results should be reliable.

DPTPP and BFB Performance Results

The DPTPP performance results were all included and found to be within the specified criteria.

Yes No _____

Comments:

The BFB performance results were all included and found to be within the specified criteria.

Yes No _____

Comments:

The (DPTPP/BFB) performance result(s) was/were reviewed and the following abundance were found to fall outside the specified criteria:

<u>Date</u>	<u>Compound</u>	<u>m/z</u>	<u>Required Abundance</u>	<u>Actual Abundance</u>
-------------	-----------------	------------	---------------------------	-------------------------

All values were within Contract limits.

All samples were extracted and analyzed within contract holding times

Yes No _____

Comments: All extraction and analysis times were within contract required holding times.

Surrogate recoveries were reviewed. The recoveries were all within the contract limits.

Yes

No. _____

Comments:

Surrogate recoveries were determined for the VOA, B/N fraction, and the acid fraction. All compounds were within the contract required limits for these fractions.

The matrix spike recovery data were reviewed. The Matrix spikes were performed and all data met contract requirements.

Yes

No. _____

Comments:

A matrix spike was performed, and the recovery of chlorobenzene was outside of the normal /normal/ limits. chlorobenzene; 172% Recovery, with limits of 60-123%. Other VOA spike recoveries were ok.

Duplicate Sample Analysis : DNA duplicate analysis was performed on this one sample but no HSL compounds were observed in either analysis. VOA duplicate analysis showed similar compounds at approximately the same concentrations, but all of the compounds observed were also in the blank analysis.

Initial calibration data were reviewed. Initial calibration data were included in the package and met all contract requirements.

Yes _____

No

Comments: The initial calibration data was not included in the raw data package, but the results from the initial calibration were included on the summary sheets so that the values were available.

Continuing calibration data were reviewed and this data met all contract requirements.

Yes _____

No

Comments: Several of the continuing calibration HSL compounds were outside of the contract limits. These compounds are listed on the enclosed sheets.

PESTICIDES

The laboratory met the pesticide linearity check criteria.

Yes _____ No _____

Comments:

N.A.

The % breakdown of 4,4DDT and of Endrin was less than 20%.

Yes _____ No _____

Comments:

N.A.

The dibutylchloroendate retention time shift was within the specified limits.

Yes _____ No _____

Comments:

N.A.

The pesticide standard compounds showed a % D of the calibration factor of no more than 15% for the quantitation runs and 20% for the confirmation runs.

Yes _____ No _____

Comments:

N.A.

BLANK ANALYSIS RESULTS

The blank analysis were reviewed. The contaminants in the blank are listed below: (*Indicates amounts above CRDL)

<u>Sample</u>	<u>Date</u>	<u>Compound</u>	<u>Conc.</u>	<u>Samples Associated with this Blank</u>
<u>VOA</u>	<u>5/6</u>	<u>Methylene chloride</u>	<u>630</u>	<u>µg/kg (all)</u>
<u> </u>	<u> </u>	<u>Acetone</u>	<u>3900</u>	<u> </u>
<u> </u>	<u> </u>	<u>2-Butanone</u>	<u>380</u>	<u> </u>
<u> </u>	<u> </u>	<u>4-methyl-2-Pentanone</u>	<u>1900</u>	<u> </u>
<u> </u>	<u> </u>	<u>Toluene</u>	<u>230</u>	<u> </u>
<u> </u>	<u> </u>	<u>Xylene</u>	<u>260</u>	<u>µg/kg (all samples)</u>

BNA All BNA Blank Compounds were T.I.C.'s,
 not H.S.L compounds.
 - No pesticide analysis. -

Remarks:

COMPOUNDS IDENTIFIED

Sample No. E.S.-SO-4 Soil Sample.

Hazardous Substances List (HSL) Compounds Detected:

Tentatively Identified Compounds Detected: SEE ATTACHED PAGE #

-Methyl Hydrocarbons-

Spectra Matching Quality

- The spectra were examined and found to be of good matching quality.
 The spectra were examined and found to be of poor matching quality due to:

Remarks:

DATA QUALIFIER DEFINITIONS
Region 8

For the purposes of this data review document the following code letters and associated definitions are provided.

- U - The material was analyzed for, but was not detected. The associated numerical value is the estimated sample quantitation limit.
- J - The associated numerical value is an estimated quantity because the amount detected is below the required limits or because quality control criteria were not met.
- UB - Estimated sample quantitation limit increased. Amount found in sample reported. Compound detected at <5 X the amount in blank (<10 X for methylene chloride, acetone, toluene and phthalates).
- UJ - Detection limit is estimated because quality control criteria were not met.
- JB - The value is an estimated amount detected below required limits and also detected in the blank.
- B - Compound was detected in the blank. Quantity reported is >5 X the amount found in the blank (>10 X for methylene chloride, acetone, toluene, and phthalates).
- R - Quality Control indicates that data is not usable (compound may or may not be present). Resampling and reanalysis is necessary for verification.
- Z - No analytical result.
- N - Presumptive evidence of presence of material (tentative identification).

DATA COMPLETENESS CHECKER

Included; no problems
 Included; problems noted in review
 Not Included

Case Narrative

Quality Control Summary Package

- Duplicate Recovery Summary (Form II)
- MS/MOS Summary (Form III)
- Reagent Blank Summary (Form IV)
- GC/MS Tuning and Mass Calibration (Form V)

Sample Data Package

- Holding Times (SMD Sample Traffic Reports)
- Organic Analysis Data Sheets (Form I; all four pages for each sample, arranged in increasing SMD number order)
- Reconstructed Ion Chromatogram(s) (RIC)
- GC/EC Chromatograms
- Quantitation Reports
- Mass Spectral Data
- EPA/NIH Mass Spectral Library Search for TIC's

Standards Data Package

- Current List of laboratory/Instrumental Detection Limits
- Initial Calibration Data (Form VI) for each instrument
- Continuing Calibration Data (Form VII) for each instrument
- Pesticide Evaluation Standards Summary (Form VIII)
- Pesticide/PCB Standards Summary (Form IX)
- Pesticide/PCB Identification (Form X; if any positive results)
- VOA and EPA Standards Reconstruction Ion Chromatograms (RIC)
- VOA and EPA Standards Quantitation Reports
- Pesticide/PCB Standard Chromatograms and Data System Printouts

Note →

Raw QC Data Package

- DPTPP and EFB mass spectra and mass listings
- Reagent Blank Data
 - Organic Analysis Data Sheets (Form I)
 - Reconstructed Ion Chromatograms (RIC)
 - Quantitation Reports
 - Mass Spectral Data
 - EPA/NIH Library Search of TIC's
 - GC/EC Chromatograms and Data System Printouts

Matrix Spike and Matrix Spike Duplicate Data ^{ONLY}

- Organic Analysis Data Sheets (Form I)
- Reconstructed Ion Chromatograms (RIC)
- Quantitation Reports
- Mass Spectral Data
- EPA/NIH Library Search of TIC's
- GC/EC Chromatograms and Data System Printouts

SAMPLE DESCRIPTION INFORMATION

for

USEPA Region VIII

<u>RMA Sample No.</u>	<u>Sample Description</u>	<u>Sample Type</u>	<u>Date Sampled</u>	<u>Date Received</u>
4909-01	ES-SO-4 Refining area	Soil	4/21/85	4/29/85
4909-02	ER-SO-1 Oily soil area	Soil	4/20/85	4/29/85
4909-03	ER-SS-1 Pond sediment	Soil	4/20/85	4/29/85

May 29, 1985

ANALYTICAL RESULTS

for

USEPA Region VIII

REFINERY VOLATILE ORGANICS

Parameter	Units	Detection Limit	4909-01 ^a		
			4909-02	4909-03	
Acetonitrile	ug/kg	300	BDL	BDL	BDL
Acrolein	ug/kg	30	BDL	BDL	BDL
Acrylonitrile*	ug/kg	-	-	-	-
Benzene	ug/kg	5	BDL	BDL	BDL
Bis(chloromethyl)ether*	ug/kg	-	-	-	-
Carbon disulfide	ug/kg	5	BDL	BDL	BDL
Chlorobenzene	ug/kg	5	BDL	BDL	BDL
Chloroform	ug/kg	5	BDL	BDL	BDL
Chloromethane	ug/kg	10	BDL	BDL	BDL
Crotonaldehyde	ug/kg	200	BDL	BDL	BDL
1,2-Dibromoethane	ug/kg	20	BDL	BDL	BDL
1,1-Dichloroethane	ug/kg	5	BDL	BDL	BDL
1,2-Dichloroethane	ug/kg	5	BDL	BDL	BDL
1,1-Dichloroethylene	ug/kg	5	BDL	BDL	BDL
1,2-Dichloroethylene	ug/kg	5	BDL	BDL	BDL
Dichloromethane	ug/kg	10	BDL	BDL	61
Dichloropropane	ug/kg	-	NA	NA	NA
1,2-Dichloropropane	ug/kg	5	BDL	BDL	BDL
1,4-Dioxane	ug/kg	150	BDL	BDL	BDL
Ethyleneimine*	ug/kg	-	-	-	-
Ethylene oxide*	ug/kg	-	-	-	-
Methyl ethyl ketone	ug/kg	10	BDL	BDL	53
Styrene	ug/kg	5	BDL	BDL	BDL
Tetrachloroethanes	ug/kg	-	NA	NA	NA
1,1,2,2-Tetrachloroethane	ug/kg	5	BDL	BDL	BDL
Tetrachloroethylene	ug/kg	5	BDL	BDL	BDL
Toluene	ug/kg	5	2000	BDL	43

BDL = Below detection limit. NA = Not analyzed. *Not recovered using Method 8240, or no analytical standard available.
 a = Detection limits 400 times higher.

ANALYTICAL RESULTS

for

USEPA Region VIII

REFINERY VOLATILE ORGANICS

Parameter	Units	Detection Limit	4909-01 ^a		
			4909-02	4909-03	
Trichloroethanes	ug/kg	-	NA	NA	NA
1,1,1-Trichloroethane	ug/kg	5	BDL	14	16
1,1,2-Trichloroethane	ug/kg	5	BDL	BDL	BDL
Trichloroethylene	ug/kg	5	BDL	BDL	BDL
Xylenes	ug/kg	-	NA	NA	NA
p-o-Xylene	ug/kg	5	BDL	BDL	BDL
m-Xylene	ug/kg	5	BDL	BDL	BDL

BDL = Below detection limit. NA = Not analyzed. *Not recovered using Method 8240, or no analytical standard available.
 a = Detection limits 400 times higher.

ANALYTICAL RESULTS

for

USEPA Region VIII

REFINERY BASE/NEUTRAL ORGANICS

Parameter	Units	Detection Limit	4909-01			4909-02			4909-03		
			4909-01	4909-02	4909-03	4909-01	4909-02	4909-03	4909-01	4909-02	4909-03
Aniline*	mg/kg	-	-	-	-	-	-	-	-	-	-
Anthracene	mg/kg	10	BDL	BDL	BDL	-	-	-	-	-	-
Benz(c)acridine*	mg/kg	-	-	-	-	-	-	-	-	-	-
Benz(a)anthracene	mg/kg	10	BDL	BDL	BDL	-	-	-	-	-	-
Benzo(b)fluoranthene	mg/kg	10	BDL	BDL	BDL	-	-	-	-	-	-
Benzo(j)fluoranthene*	mg/kg	-	-	-	-	-	-	-	-	-	-
Benzo(k)fluoranthene	mg/kg	10	BDL	BDL	BDL	-	-	-	-	-	-
Benzo(a)pyrene	mg/kg	10	BDL	BDL	BDL	-	-	-	-	-	-
Benzyl chloride*	mg/kg	-	-	-	-	-	-	-	-	-	-
Bis(2-ethylhexyl)phthalate	mg/kg	10	BDL	24	11	-	-	-	-	-	-
Butyl benzyl phthalate	mg/kg	5	BDL	BDL	BDL	-	-	-	-	-	-
2-Chloronaphthalene	mg/kg	5	BDL	BDL	BDL	-	-	-	-	-	-
Chrysene	mg/kg	10	BDL	BDL	BDL	-	-	-	-	-	-
Dibenz(a,h)acridine*	mg/kg	-	-	-	-	-	-	-	-	-	-
Dibenz(a,j)acridine	mg/kg	10	BDL	BDL	BDL	-	-	-	-	-	-
Dibenz(a,h)anthracene	mg/kg	10	BDL	BDL	BDL	-	-	-	-	-	-
Dibenzo(a,e)pyrene*	mg/kg	-	-	-	-	-	-	-	-	-	-
Dibenzo(a,h)pyrene*	mg/kg	-	-	-	-	-	-	-	-	-	-
Dibenzo(a,i)pyrene*	mg/kg	-	-	-	-	-	-	-	-	-	-
Di-n-butyl phthalate	mg/kg	5	BDL	BDL	BDL	-	-	-	-	-	-
Dichlorobenzenes	mg/kg	-	NA	NA	NA	-	-	-	-	-	-
o-Dichlorobenzene	mg/kg	5	BDL	BDL	BDL	-	-	-	-	-	-
m-Dichlorobenzene	mg/kg	5	BDL	BDL	BDL	-	-	-	-	-	-
p-Dichlorobenzene	mg/kg	5	BDL	BDL	BDL	-	-	-	-	-	-
Diethyl phthalate	mg/kg	50	BDL	BDL	BDL	-	-	-	-	-	-
7,12-Dimethylbenz(a)anthracene	mg/kg	5	BDL	BDL	BDL	-	-	-	-	-	-
Dimethyl phthalate	mg/kg	20	BDL	BDL	BDL	-	-	-	-	-	-
2,4-Dinitrotoluene	mg/kg	10	BDL	BDL	BDL	-	-	-	-	-	-
Di-n-octyl phthalate	mg/kg	5	BDL	BDL	BDL	-	-	-	-	-	-

BDL = Below detection limit. NA = Not analyzed. *Not recovered using Method 8270, or no analytical standard available.

ANALYTICAL RESULTS

for

USEPA Region VIII

REFINERY BASE/NEUTRAL ORGANICS (Cont.)

Parameter	Units	Detection Limit	4909-01			4909-02			4909-03		
			4909-01	4909-02	4909-03	4909-01	4909-02	4909-03	4909-01	4909-02	4909-03
Fluoranthene	mg/kg	5	BDL								
Indene	mg/kg	10	BDL								
Indeno(1,2,3-c,d)pyrene	mg/kg	15	BDL								
Methylbenz(c)phenanthrene	mg/kg	5	BDL								
3-Methylicholanthrene	mg/kg	10	BDL								
Methyl chrysene*	mg/kg	-	-	-	-	-	-	-	-	-	-
1-Methylnaphthalene	mg/kg	10	BDL								
Naphthalene	mg/kg	5	BDL								
Naphthylamine	mg/kg	-	NA								
1-Naphthylamine	mg/kg	100	BDL								
2-Naphthylamine	mg/kg	100	BDL								
5-Nitroacenaphthene*	mg/kg	-	-	-	-	-	-	-	-	-	-
p-Nitroaniline	mg/kg	100	BDL								
Nitrobenzene	mg/kg	10	BDL								
Phenanthrene	mg/kg	5	BDL								
Pyrene	mg/kg	10	BDL								
Pyridine*	mg/kg	-	-	-	-	-	-	-	-	-	-
Quinoline	mg/kg	100	BDL								
Trichlorobenzenes	mg/kg	-	NA								
1,2,4-Trichlorobenzene	mg/kg	5	BDL								
Trimethylbenz(a)anthracene*	mg/kg	-	-	-	-	-	-	-	-	-	-

BDL = Below detection limit. NA = Not analyzed. *Not recovered using Method 8270, or no analytical standard available.

ANALYTICAL RESULTS

for

USEPA Region VIII

REFINERY ACID ORGANICS

Parameter	Units	Detection Limit	4909-01 4909-02 4909-03		
			4909-01	4909-02	4909-03
Benzene thiol*	mg/kg	-	-	-	-
p-Chloro-m-cresol	mg/kg	5	BDL	BDL	BDL
2-Chlorophenol	mg/kg	5	BDL	BDL	BDL
Cresols	mg/kg	-	NA	NA	NA
o-Cresol	mg/kg	5	BDL	BDL	BDL
p & m-Cresol	mg/kg	5	BDL	BDL	BDL
2,4-Dimethylphenol	mg/kg	5	BDL	BDL	BDL
4,6-Dinitro-o-cresol	mg/kg	5	BDL	BDL	BDL
2,4-Dinitrophenol	mg/kg	200	BDL	BDL	BDL
4-Nitrophenol	mg/kg	20	BDL	BDL	BDL
Pentachlorophenol	mg/kg	5	BDL	BDL	BDL
Phenol	mg/kg	10	BDL	BDL	BDL
Trichlorophenols	mg/kg	-	NA	NA	NA
2,4,5-Trichlorophenol	mg/kg	5	BDL	BDL	BDL
2,4,6-Trichlorophenol	mg/kg	5	BDL	BDL	BDL

BDL = Below detection limit. NA = Not analyzed. *Not recovered using Method 8270, or no analytical standard available.

ANALYTICAL RESULTS

for

USEPA Region VIII

REFINERY ACID ORGANICS

Parameter	Units	Detection Limit	4909-01			4909-02			4909-03		
			4909-01	4909-02	4909-03	4909-01	4909-02	4909-03	4909-01	4909-02	4909-03
Benzene-thiol*	mg/kg	-	-	-	-	-	-	-	-	-	-
p-Chloro-m-cresol	mg/kg	5	BDL								
2-Chlorophenol	mg/kg	5	BDL								
Cresols	mg/kg	-	NA								
o-Cresol	mg/kg	5	BDL								
p & m-Cresol	mg/kg	5	BDL								
2,4-Dimethylphenol	mg/kg	5	BDL								
4,6-Dinitro-o-cresol	mg/kg	5	BDL								
2,4-Dinitrophenol	mg/kg	200	BDL								
4-Nitrophenol	mg/kg	20	BDL								
Pentachlorophenol	mg/kg	5	BDL								
Phenol	mg/kg	10	BDL								
Trichlorophenols	mg/kg	-	NA								
2,4,5-Trichlorophenol	mg/kg	5	BDL								
2,4,6-Trichlorophenol	mg/kg	5	BDL								

BDL = Below detection limit. NA = Not analyzed. *Not recovered using Method 8270, or no analytical standard available.

Sample Number

H47-H1848

4909-01
100ml

JULY 31

Organics Analysis Data Sheet
(Page 1)

Laboratory Name: RNAL
Lab Sample ID No: 4909-01
Sample Matrix: STO soil
Data Release Authorized By: _____

Case No: _____
QC Report No: _____
Contract No: _____
Date Sample Received: _____

$$\frac{\text{ug}}{\text{kg}} \times \frac{0.005}{\text{ug}} \times \frac{0.020}{\text{ug}} = \frac{0.00246}{\text{kg}} \text{ kg.}$$

$$= 406 \text{ ug/kg}$$

Volatile Compounds

Concentration: Low Medium (Circle One)
Date Extracted/Prepared: _____
Date Analyzed: 5/6/85
Conc/Dil Factor: x406 pH _____

Percent Moisture: _____
Percent Moisture (Decanted): _____

CAS Number		ug/l or ug/Kg (Circle One)
74-87-3	Chloromethane	
74-83-9	Bromomethane	
75-01-4	Vinyl Chloride	
75-00-3	Chloroethane	
75-09-2	Methylene Chloride	<u>2500</u> <i>SB</i>
67-64-1	Acetone	<u>29600</u> <i>BR</i>
75-15-0	Carbon Disulfide	
75-35-4	1, 1-Dichloroethene	
75-34-3	1, 1-Dichloroethane	
156-60-5	Trans-1, 2-Dichloroethene	
67-66-3	Chloroform	
107-06-2	1, 2-Dichloroethane	
78-93-3	2-Butanone	2000 <i>SB</i>
71-55-6	1, 1, 1-Trichloroethane	
56-23-5	Carbon Tetrachloride	
108-05-4	Vinyl Acetate	
75-27-4	Bromodichloromethane	

CAS Number	ug/l or ug/Kg (Circle One)
79-34-5	1, 1, 2, 2-Tetrachloroethane
78-87-5	1, 2-Dichloropropane
10061-02-6	Trans-1, 3-Dichloropropene
79-01-6	Trichloroethene
124-48-1	Dibromochloromethane
79-00-5	1, 1, 2-Trichloroethane
71-43-2	Benzene
10061-01-5	cis-1, 3-Dichloropropene
110-75-8	2-Chloroethylvinylether
75-25-2	Bromoform
591-78-6	2-Hexanone
108-10-1	4-Methyl-2-Pentanone
127-18-4	Tetrachloroethene
108-88-3	Toluene
108-90-7	Chlorobenzene
100-41-4	Ethylbenzene
100-42-5	Styrene
	Total Xylenes

Data Reporting Qualifiers

For reporting results to EPA, the following results qualifiers are used.

Additional flags or footnotes explaining results are encouraged. However, the definition of each flag must be explicit.

Value If the result is a value greater than or equal to the detection limit, report the value

C This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides $\geq 10 \text{ ng}/\text{ul}$ in the final extract should be confirmed by GC/MS

U Indicates compound was analyzed for but not detected. Report the minimum detection limit for the sample with the U (e.g., 10U) based on necessary concentration/dilution actions. (This is not necessarily the instrument detection limit.) The footnote should read: U-Compound was analyzed for but not detected. The number is the minimum attainable detection limit for the sample.

B This flag is used when the analyte is found in the blank as well as a sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action.

J Indicates an estimated value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicates the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero (e.g., 10J).

Other Other specific flags and footnotes may be required to properly define the results. If used, they must be fully described and such description attached to the data summary report.

Sample Number
4909-01
100 uL

Organics Analysis Data Sheet
(Page 4)

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or ug/kg)
1.	CG Hydrocarbon	VOA	267	1600
2.	Unknown Hydrocarbon	VOA	321	6400
3.	CG Hydrocarbon	VOA	360	6400
4.	Unknown	VOA	393	2500
5.	Hexane / Octene	VOA	470	3600
6.	Unknown Hydrocarbon	VOA	499	5100
7.	Trimethyl-Cyclohexane	VOA	512	31000
8.	C9 Unsaturated Hydrocarbon	VOA	527	5500
9.	Trimethyl-Cyclohexane	VOA	538	6300
10.	C9 Unsaturated Hydrocarbon	VOA	567	14000
11.	C9 Unsaturated Hydrocarbon	VOA	582	23000
12.	Dimethyl-Cyclohexane	VOA	600	15000
13.	Unknown	VOA	610	20000
14.	Unknown	VOA	635	5600
15.	C10 Hydrocarbon	VOA	673	9200
16.				
17.				
18.				
19.				
20.				
21.				
22.				
23.				
24.				
25.				
26.				
27.				
28.				
29.				
30.				

Sample Number

4909-01

DUPLICATE

Organics Analysis Data Sheet
(Page 1)

Laboratory Name: Pac Mountain Polymer Lab

Lab Sample ID No: S4909A133

Sample Matrix: SOIL - GLYME

Data Release Authorized By: _____

Case No: _____

QC Report No: _____

Contract No: _____

Date Sample Received: _____

Volatile Compounds

Concentration: Low Medium (Circle One)

Date Extracted/Prepared: -

Date Analyzed: 5/7/85

Conc/Dil Factor: 395 kg⁻¹ pH -

Percent Moisture: -

Percent Moisture (Decanted): -

$$\left(\frac{0.005\text{L}}{1.00253\text{kg}} \right) \left(\frac{20\text{ml}}{0.100\text{ml}} \right) = 395 \text{ L kg}^{-1}$$

CAS Number		ug/l or ug/Kg (Circle One)	CAS Number	ug/l or ug/Kg (Circle One)
74-87-3	Chloromethane		79-34-5	1, 1, 2-Tetrachloroethane
74-83-9	Bromomethane		78-87-5	1, 2-Dichloropropane
75-01-4	Vinyl Chloride		10061-02-6	Trans-1, 3-Dichloropropene
75-00-3	Chloroethane		79-01-6	Trichloroethene
75-09-2	Methylene Chloride	2 J ✓	124-48-1	Dibromochloromethane
67-64-1	Acetone	28 ✓	79-00-5	1, 1, 2-Trichloroethane
75-15-0	Carbon Disulfide		71-43-2	Benzene
75-35-4	1, 1-Dichloroethene		10061-01-5	cis-1, 3-Dichloropropene
75-34-3	1, 1-Dichloroethane		110-75-8	2-Chloroethylvinylether
156-60-5	Trans-1, 2-Dichloroethene		75-25-2	Bromoform
67-66-3	Chloroform		591-78-6	2-Hexanone
107-06-2	1, 2-Dichloroethane		108-10-1	4-Methyl-2-Pentanone
78-93-3	2-Butanone	2 NO ✓	127-18-4	Tetrachloroethene
71-55-6	1, 1, 1-Trichloroethane	1 J ✓	108-88-3	Toluene
56-23-5	Carbon Tetrachloride		108-90-7	Chlorobenzene
108-05-4	Vinyl Acetate		100-41-4	Ethylbenzene
75-27-4	Bromodichloromethane		100-42-5	Styrene
<u>1,4-DIOXANE</u>		<u>39400</u>	Total Xylenes	

Data Reporting Qualifiers

For reporting results to EPA, the following results qualifiers are used

Additional flags or footnotes explaining results are encouraged. However, the definition of each flag must be explicit

V Value If the result is a value greater than or equal to the detection limit, report the value

C This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides $\geq 10 \text{ ng}/\text{ul}$ in the final extract should be confirmed by GC/MS

U Indicates compound was analyzed for but not detected. Report the minimum detection limit for the sample with the U (e.g., 10U) based on necessary concentration/dilution factors. (This is not necessarily the instrument detection limit.) The footnote should read U. Compound was analyzed for but not detected. The number is the minimum attainable detection limit for the sample

B This flag is used when the analyte is found in the blank as well as a sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action

J Indicates an estimated value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicates the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but

Other Other specific flags and footnotes may be required to properly define the results. If used, they must be fully described and such description attached to the data summary report

Sample Number
4909-01

DUPPLICATE

Organics Analysis Data Sheet
(Page 4)

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or ug/kg)
1.	C ₆ ALIPHATIC	VOA	360	1.6
2.	UNKNOWN	VOA	471	18
3.	C ₈₋₉ ALIPHATIC	VOA	500	1.9
4.3073-66-3	1,1,3-TRIMETHYL CYCLOHEXANE	VOA	513	25
5.	C ₃ -SUBSTITUTED CYCLOHEXANE	VOA	527	4.4
6.	C ₃ -SUBSTITUTED CYCLOHEXANE	VOA	539	4.0
7.	C ₃ -SUBSTITUTED CYCLOHEXANE	VOA	568	6.4
8.	- C ₉ - ALIPHATIC	VOA	583	9.2
9.	C ₁₀ - ALIPHATIC	VOA	611	18
10.	UNKNOWN	VOA	636	3.3
11.	C ₁₀ - ALIPHATIC	VOA	674	4.7
12.				
13.				
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29.				
30.				

Environmental Protection Agency, CLP Sample Management Office,
P. O. Box 818, Alexandria, Virginia 22313 703/657-2490

Sample Number

4904-01

Acid

Organics Analysis Data Sheet
(Page 2)

Semivolatile Compounds

Concentration: Low Medium (Circle One)

Date Extracted/Prepared: 5-9-86

Date Analyzed: 5-19-85

Conc/Dil Factor: 0.100 5.0

$$0.001 \text{ mg/kg} \times \frac{1 \text{ mg}}{1000 \text{ ug}} = 4.13$$

J. J. Kelly

mg/kg

CAS Number		<i>mg/kg</i> ND
		(Circle One)
62-75-9	N-Nitrosodimethylamine	<i>ND</i>
108-95-2	Phenol	
62-53-3	Aniline	
111-44-4	bis(2-Chloroethyl)Ether	
95-57-8	2-Chlorophenol	
541-73-1	1, 3-Dichlorobenzene	
106-46-7	1, 4-Dichlorobenzene	
100-51-6	Benzyl Alcohol	
95-50-1	1, 2-Dichlorobenzene	
95-48-7	2-Methylphenol	
39638-32-9	bis(2-chloroisopropyl)Ether	
106-44-5	4-Methylphenol	
621-64-7	N-Nitroso-Di-n-Propylamine	
67-72-1	Hexachloroethane	
98-95-3	Nitrobenzene	
78-59-1	Isophorone	
88-75-5	2-Nitrophenol	
105-67-9	2, 4-Dimethylphenol	
65-85-0	Benzoic Acid	
111-91-1	bis(2-Chloroethoxy)Methane	
120-83-2	2, 4-Dichlorophenol	
120-82-1	1, 2, 4-Trichlorobenzene	
91-20-3	Naphthalene	
106-47-8	4-Chloroaniline	
87-68-3	Hexachlorobutadiene	
59-50-7	4-Chloro-3-Methylphenol	
91-57-6	2-Methylnaphthalene	
77-47-4	Hexachlorocyclopentadiene	
88-06-2	2, 4, 6-Trichlorophenol	
95-95-4	2, 4, 5-Trichlorophenol	
91-58-7	2-Chloronaphthalene	
88-74-4	2-Nitroaniline	
131-11-3	Dimethyl Phthalate	
208-96-8	Arenaphthylene	
99-09-2	3-Nitroaniline	

CAS Number		<i>mg/kg</i> ND
		(Circle One)
83-32-9	Acenaphthene	<i>ND</i>
51-28-5	2, 4-Dinitrophenol	
100-02-7	4-Nitrophenol	
132-64-9	Dibenzofuran	
121-14-2	2, 4-Dinitrotoluene	
606-20-2	2, 6-Dinitrotoluene	
84-66-2	Diethylphthalate	
7005-72-3	4-Chlorophenyl-phenylether	
86-73-7	Fluorene	
100-01-6	4-Nitroaniline	
534-52-1	4, 6-Dinitro-2-Methylphenol	
86-30-6	N-Nitrosodiphenylamine (1)	
101-55-3	4-Bromophenyl-phenylether	
118-74-1	Hexachlorobenzene	
87-86-5	Pentachlorophenol	
85-01-8	Phenanthrene	
120-12-7	Anthracene	
84-74-2	Di-n-Butylphthalate	
206-44-0	Fluoranthene	
92-87-5	Benzidine	
129-00-0	Pyrene	
85-68-7	Butylbenzylphthalate	
91-94-1	3, 3'-Dichlorobenzidine	
56-55-3	Benz(a)Anthracene	
117-81-7	bis(2-Ethylhexyl)Phthalate	
218-01-9	Chrysene	
117-84-0	Di-n-Octyl Phthalate	
205-99-2	Benzo(b)Fluoranthene	
207-08-9	Benzo(k)Fluoranthene	
50-32-8	Benzo(a)Pyrene	
193-39-5	Indeno(1, 2, 3-cd)Pyrene	
53-70-3	Dibenzo(a, h)Anthracene	
191-24-2	Benzo(a, h, i)Perylene	

(1) Cannot be separated from diphenylamine.

Sample Number

4907-01

B/N

Organics Analysis Data Sheet
(Page 2)

Semi-volatile Compounds

Concentration: Low Medium (Circle One)

Date Extracted/Prepared: 5-9-85

Date Analyzed: 5-10-85

Conc/Dil Factor: Half 5.0 mL

$$\frac{1}{400 \text{ mL}} \cdot \frac{1}{1000} = 4/3 \text{ (mg/kg)}$$

CAS Number	Description	mg/kg
		Detectable? (Circle One)
62-75-9	N-Nitrosodimethylamine	<u>yes</u>
108-95-2	Phenol	
62-53-3	Aniline	
111-44-4	bis(2-Chloroethyl)Ether	<u>yes</u>
95-57-8	2-Chlorophenol	<u>yes</u>
541-73-1	1, 3-Dichlorobenzene	
106-46-7	1, 4-Dichlorobenzene	
100-51-8	Benzyl Alcohol	
95-50-1	1, 2-Dichlorobenzene	
95-48-7	2-Methoxyphenol	
39638-32-9	bis(2-chloroisopropyl)Ether	
106-44-5	4-Methoxyphenol	
621-64-7	N-Nitroso-Di-n-Propylamine	
67-72-1	Hexachloroethane	
98-95-3	Nitrobenzene	
78-59-1	Isophorone	
88-75-5	2-Nitrophenol	
105-67-9	2, 4-Dimethoxyphenol	
65-85-0	Benzoic Acid	
111-91-1	bis(2-Chloroethoxy)Methane	
120-83-2	2, 4-Dichlorophenol	
120-82-1	1, 2, 4-Trichlorobenzene	
91-20-3	Naphthalene	
106-47-8	4-Chloroaniline	
87-68-3	Hexachlorobutadiene	
59-50-7	4-Chloro-3-Methoxyphenol	
91-57-6	2-Methylnaphthalene	
77-47-4	Hexachlorocyclopentadiene	
88-06-2	2, 4, 6-Trichlorophenol	
95-95-4	2, 4, 5-Trichlorophenol	
91-58-7	2-Chloronaphthalene	
88-74-4	2-Nitroaniline	
131-11-3	Dimethyl Phthalate	
208-96-8	Acenaphthylene	
99-09-2	3-Nitroaniline	

CAS Number	Description	Detectable? (Circle One)
83-32-9	Acenaphthene	<u>yes</u>
51-28-5	2, 4-Dinitrophenol	<u>yes</u>
100-02-7	4-Nitrophenol	
132-64-9	Dibenzofuran	
121-14-2	2, 4-Dinitrotoluene	
605-20-2	2, 6-Dinitrotoluene	
64-66-2	Diethylphthalate	
7005-72-3	4-Chlorophenyl-phenylether	
86-73-7	Fluorene	
100-01-6	4-Nitroaniline	
534-52-1	4, 6-Dinitro-2-Methoxyphenol	
86-30-6	N-Nitrosodiphenylamine (1)	
101-55-3	4-Bromophenyl-phenylether	
118-74-1	Hexachlorobenzene	
87-86-5	Pentachlorophenol	
85-01-8	Phenanthrene	
120-12-7	Anthracene	
84-74-2	Di-n-Butylphthalate	
206-44-0	Fluoranthene	
92-87-5	Benzidine	
129-00-0	Pyrene	
85-68-7	Butylbenzylphthalate	
91-94-1	3, 3'-Dichlorobenzidine	
56-55-3	Benz(a)Anthracene	
117-81-7	bis(2-Ethylhexyl)Phthalate	
218-01-9	Chrysene	
117-84-0	Di-n-Octyl Phthalate	
205-99-2	Benz(b)Fluoranthene	
207-08-9	Benz(k)Fluoranthene	
50-32-8	Benz(a)Pyrene	
193-39-5	Indeno(1, 2, 3-cd)Pyrene	
53-70-3	Di-benz(a, h)Anthracene	
191-24-2	Benzog. h, i)Perylene	

(1)-Cannot be separated from diphenylamine

Sample Number

4909-01

Organics Analysis Data Sheet
(Page 2)

Acid Duplicate

V-7167

Semivolatile Compounds

Concentration: Low Medium (Circle One)

Date Extracted/Prepared: 5-9-85

Date Analyzed: 5-10-85

Conc/Dil Factor: 1000 x 5.0ml

$$0.001396 \times \frac{1}{1000} = 3.97 \times 10^{-6}$$

100/47

34 mg/kg

mylk

~~mg/kg/Kg~~
(Circle One)

CAS
Number

62-75-9	N-Nitrosodimethylamine	<input checked="" type="radio"/>
108-95-2	Phenol	
62-53-3	Aniline	
111-44-4	bis(2-Chloroethyl)Ether	
95-57-8	2-Chlorophenol	
541-73-1	1, 3-Dichlorobenzene	
106-46-7	1, 4-Dichlorobenzene	
100-51-6	Benzyl Alcohol	
95-50-1	1, 2-Dichlorobenzene	
95-48-7	2-Methylphenol	
39638-32-9	bis(2-chloroisopropyl)Ether	
106-44-5	4-Methylphenol	
621-64-7	N-Nitroso-Di-n-Propylamine	
67-72-1	Hexachloroethane	
98-95-3	Nitrobenzene	
78-59-1	Isophorone	
88-75-5	2-Nitrophenol	
105-67-9	2, 4-Dimethylphenol	
65-85-0	Benzoic Acid	
111-91-1	bis(2-Chloroethoxy)Methane	
120-83-2	2, 4-Dichlorophenol	
120-82-1	1, 2, 4-Trichlorobenzene	
91-20-3	Naphthalene	
106-47-8	4-Chloroaniline	
87-68-3	Hexachlorobutadiene	
59-50-7	4-Chloro-3-Methylphenol	
91-57-6	2-Methylnaphthalene	
77-47-4	Hexachlorocyclopentadiene	
88-06-2	2, 4, 6-Trichlorophenol	
95-95-4	2, 4, 5-Trichlorophenol	
91-58-7	2-Chloronaphthalene	
88-74-4	2-Nitroaniline	
131-11-3	Dimethyl Phthalate	
208-96-8	Acenaphthylene	
99-09-2	3-Nitroaniline	

CAS Number	Concentration (Circle One)
83-32-9	Acenaphthene
51-28-5	2, 4-Dinitrophenol
100-02-7	4-Nitrophenol
132-64-9	Dibenzofuran
121-14-2	2, 4-Dinitrotoluene
606-20-2	2, 6-Dinitrotoluene
84-66-2	Diethylphthalate
7005-72-3	4-Chlorophenyl-phenylether
86-73-7	Fluorene
100-01-6	4-Nitroaniline
534-52-1	4, 6-Dinitro-2-Methylphenol
86-30-6	N-Nitrosodiphenylamine (1)
101-55-3	4-Bromophenyl-phenylether
118-74-1	Hexachlorobenzene
87-86-5	Pentachlorophenol
85-01-8	Phenanthrene
120-12-7	Anthracene
84-74-2	Di-n-Butylphthalate
206-44-0	Fluoranthene
92-87-5	Benzidine
129-00-0	Pyrene
85-68-7	Butylbenzylphthalate
91-94-1	3, 3'-Dichlorobenzidine
56-55-3	Benz(a)Anthracene
117-81-7	bis(2-Ethylhexyl)Phthalate
218-01-9	Chrysene
117-84-0	Di-n-Octyl Phthalate
205-99-2	Benz(b)Fluoranthene
207-08-9	Benz(k)Fluoranthene
50-32-8	Benz(a)Pyrene
193-39-5	Indeno(1, 2, 3-cd)Pyrene
53-70-3	Dibenz(a, h)Anthracene
191-24-2	Benz(o, h,i)Perylene

(*) Cannot be separated from diphenylamine

Sample Number
4909-01

B/N

Duplicate

Organics Analysis Data Sheet
(Page 2)

Semi-volatile Compounds

Concentration: Low Medium (Circle One)

Date Extracted/Prepared: 5-8-85

Date Analyzed: 5-21-85

Conc/Dil Factor: 1/2 5.0ml
~~ml~~ ~~600/47 1000~~ ~~3.4ml~~

2001/09

10/17/85

CAS Number		(Circle One)
62-75-9	N-Nitrosodimethylamine	ND
108-95-2	Pheno!	
62-53-3	Aniline	
111-44-4	bis(2-Chloroethyl)Ether	
95-57-8	2-Chlorophenol	
541-73-1	1, 3-Dichlorobenzene	
106-46-7	1, 4-Dichlorobenzene	
100-51-6	Benzyl Alcohol	
95-50-1	1, 2-Dichlorobenzene	
95-48-7	2-Methylphenol	
39638-32-9	bis(2-chloroisopropyl)Ether	
106-44-5	4-Methylphenol	
621-64-7	N-Nitroso-Di-n-Propylamine	
67-72-1	Hexachlorethane	
98-95-3	Nitrobenzene	
78-59-1	Iscophorone	
88-75-5	2-Nitrophenol	
105-67-5	2, 4-Dimethylphenol	
65-05-0	Benzoic Acid	
111-91-1	bis(2-Chloroethyl)Methane	
120-03-2	2, 4-Dichlorophenol	
120-02-1	1, 2, 4-Trichlorobenzene	
91-20-3	Naphthalene	
106-47-8	4-Chloroaniline	
87-08-3	Hexachlorobutadiene	
59-59-7	4-Chloro-3-Methylphenol	
91-57-6	2-Methylnaphthalene	
77-47-4	Hexachlorocyclopentadiene	
53-06-2	2, 4, 6-Trichlorophenol	
95-95-4	2, 4, 5-Trichlorophenol	
91-59-7	2-Chloronaphthalene	
108-74-4	2-Nitroaniline	
131-11-3	Dimethyl Phthalate	
220-96-6	A-epoxybutane	
99-09-2	3-Nitroaniline	

CAS Number		(Circle One)
83-32-9	Acenaphthene	ND
51-23-5	2, 4-Dinitrophenol	
100-02-7	4-Nitrophenol	
132-64-9	Dibenzofuran	
121-14-2	2, 4-Dinitrotoluene	
603-20-2	2, 6-Dinitrotoluene	
34-66-2	Diethylphthalate	
7005-72-3	4-Chlorophenyl-phenylether	
86-73-7	Fluorene	
100-01-6	4-Nitroaniline	
534-62-1	4, 6-Dinitro-2-Methylphenol	
86-30-6	N-Nitrodiphenylamine (I)	
101-55-3	4-Bromophenyl-phenylether	
118-74-1	Hexachlorobenzene	
97-06-5	Pentachlorophenol	
85-01-8	Phenanthrene	
120-12-7	Anthracene	
84-74-2	Di-n-Butylphthalate	
206-44-0	Fluoranthene	
92-87-5	Benzidine	
129-00-0	Pyrene	
35-61-7	Butyl(Benzyl)phthalate	
31-94-1	3, 3'-Dichlorobenzidine	
56-55-3	Benz(a)Anthracene	
117-01-7	bis(2-Ethylhexyl)Phthalate	
218-01-0	Chrysene	
117-84-0	Di-n-Octyl Phthalate	
203-00-2	Benz(b)Fluoranthene	
207-08-9	Benz(k)Fluoranthene	
50-32-4	(Benzol)Pyrene	
183-30-5	Indundol-1, 2, 3-ediprene	
43-70-3	Dihydro-1, 6-Anthracyne	
101-24-2	Benz(a, h)Perylene	

(I) Cannot be separated from N-nitrosodiamine

REGION VIII SUMMARY OF DATA QUALITY ASSURANCE REVIEW

Case No. 4242 Project No. 4242

Site Empire State Refinery

Contractor Laboratory _____

Data Reviewer Randy Greaves Date of Review 10/4/85

Sample Matrix water and soil

Sample No. H1839 H1843(s) H1849(s)

H1839 H1844(s)

H1840 H1845(s)

H1841 H1846(s)

H1842 H1847(s)

Data are acceptable for use

Data are acceptable for use with qualification noted above

Data are preliminary - pending action or verification

Data are unacceptable

Action required by DPO?

No Yes Following items require action

Action required by Project Officer (PO)?

No Yes _____

Following are our findings:

① holding times were exceeded for samples H1839 and H1842, because samples had to be re-analyzed.

② 1,2-Dichloropropane was the only CCC compound with a RF < 0.30 for the initial calibration. None of the SCC's were out of specification for the initial calibration. SPCC's and CCL's were all within contract requirements for the continuing calibration. Several of the Non-SPCC's + CCC's compounds were not within contract limits and are listed on the calibration review sheets.

③ Methylene Chloride, acetone and butanone were the most common blank VOA compounds. Phthalate was the most common extractable contaminant. Several VOA, HSL compounds were indicated as present at below the CRDL but because of the low concentrations these compounds are not significant. Qualification were made to the organic analysis sheets based on the blank observations. These qualifications are listed on the HSL compound detected sheets.

④ Some of the advisory limits on the matrix spike and matrix spike duplicate were exceeded and the particular compounds are listed under the MS/MSD review section. No action was taken as a result of the MS/MSD results.

DPTPP and BFB Performance Results

The DPTPP performance results were all included and found to be within the specified criteria.

Yes No

Comments: All results good.

The BFB performance results were all included and found to be within the specified criteria.

Yes No

Comments: All raw data and summary sheet results were within contract requirement.

The (DPTPP/BFB) performance result(s) was/were reviewed and the following abundance were found to fall outside the specified criteria:

<u>Date</u>	<u>Compound</u>	<u>m/z</u>	<u>Required Abundance</u>	<u>Actual Abundance</u>
<u>No results</u>				

All samples were extracted and analyzed within contract holding times

Yes No

Comments: Holding times were exceeded for extraction of samples H1839 and H1842 because these had to be re-analyzed because surrogate recoveries were out of control. This was an unavoidable situation and re-analysis results with exceeded holding times and good surrogate recoveries are preferable to poor surrogate recoveries and good holding times.

Initial calibration data were reviewed. Initial calibration data were included in the package and met all contract requirements.

Yes

No ✓ (because of 1,2-Dichloropropane)

Comments:	<u>Compound</u>	<u>RF</u>	<u>Spec.</u>	
March 25,	2-Butanone	0.028	i.e., < 0.05	— Flag R neg. - Flag J pos.
	Vinyl acetate	0.007	" "	"
	2-Chloroethylvinylether	0.04	i.e., < 0.05	"

1,2-Dichloropropane. $\bar{R}F = 0.268$ i.e., < 0.30 for Spec CCP Flag J or UJ

<u>Compound</u>	<u>% RSD</u>	<u>Spec</u>	
Acetone	44.8	i.e., > 30.5%	
Vinyl acetate	39.4	i.e., > 30.5%	Flag UJ for neg. results.
4-chloroaniline	41.3	i.e., > 30.5%	Flag J for pos. results.
3-Nitroaniline	58.9	i.e., > 30.5%	

The above values were flagged as indicated above. Only one of the above compounds was a Spec or an CCP i.e. - 1,2-Dichloropropane. This compound was flagged J if present and UJ if not present. All other contract requirements were met. Continuing calibration data were reviewed and this data met all contract requirements.

Yes ✓

No

Comments: The compounds given below were made as indicated. None of the compounds given below were Spec or CCP.

<u>Compound</u>	<u>RF</u>	<u>Spec</u>
Vinyl acetate	0.014	< 0.05
2-chloroethylvinylether	0.011	< 0.05
2-Butanone	0.048	< 0.05

① pos. Flag J ↑

② Neg Flag R

↑
(VOA's)
ONLY

<u>Compound</u>	<u>% D</u>	<u>Spec.</u>	<u>Date</u>
Methylchloride	-34	25	Apr 26 4/27
acetone	-49, 63	25	"
2-butanone	-39, 82		
1,1,1-trichloroethane	42, 59		
Carbon Tetrachloride	29, 56		
Vinyl chloride	100, 100		
trans-1,3-dichloropropene	26		
cis-1,3-dichloropropene	26		
2-Hexanone	30		
1,1,2,2-Tetrachloroethane	41, 57	25	4/16 4/17
4-Methyl-2-Pentanone	24		4/27

↑ pos. Flag J.
(VOA's ONLY) Neg Flag UJ

Initial calibration data were reviewed. Initial calibration data were included in the package and met all contract requirements.

Yes _____

No _____

Comments: see last page.

Continuing calibration data were reviewed and this data met all contract requirements.

Yes

No _____

None of the compounds below were seen at spec!
Comments: continuation of last page. Below are all TNA calibration results.

<u>Compound</u>	<u>% D</u>	<u>Spec.</u>
Anthracene.	27%	25
bis(2-chloroethyl) Ether	60%	25
4-Chloroaniline	45%	25
3-Nitroaniline	69%	25
Benzoic acid	47+27%	25
4-Naphthylene	34	25
4-Nitroaniline.	47	25
Benzo[k] fluoranthene	26	25
Aniline,	51	25
2,4-dinitrophenol	29	25
4-Nitrophenol.	37	25

<u>Compound</u>	<u>% D</u>	<u>RF</u>	<u>Spec</u>	<u>pos. Flag J</u>
Benzidine	-	-	0.05	pos. Flag J
2,3-dichlorobenzidine	-	-	0.05	neg. Flag R

3-Nitroaniline. 0.034 > 0.05

<u>Compound</u>	<u>% D</u>	<u>Spec.</u>
4-Bromophenyl phenyl ether	34	25
1,4-dichlorobenzene	43	25
bis (2-Ethylhexyl) phthalate	28	25



pos. Flag J

neg. Flag UJ

Surrogate recoveries were reviewed. The recoveries were all within the contract limits.

Yes No

Comments: All values reported on the water and soil surrogate summary report forms were within contract limits. Individual calculations of compound recoveries show that summary sheet values are correct.

The matrix spike recovery data were reviewed. The Matrix spikes were performed and all data met contract requirements.

Yes No

Comments: Phenol + [] both were outside the QC advisory range for Low Water: 4-Nitrophenol [] % rec. from a water spike.

All water RPD were ok.

Low level soil: 1,2,4-Trichlorobenzene RPD was to high.
1,4-Dichlorobenzene RPD was also to high.

Medium level soil: 2,4-Dinitrotoluene — % rec out of control.
Phenol — % rec. out of control.

All RPD's were within spec. for the medium level soil.

NA = Not Applicable

PESTICIDES

The laboratory met the pesticide linearity check criteria.

Yes _____ No _____

Comments:

NA

The % breakdown of 4,4DDT and of Endrin was less than 20%.

Yes _____ No _____

Comments:

NA

The dibutylchloroendate retention time shift was within the specified limits.

Yes _____ No _____

Comments:

NA

The pesticide standard compounds showed a % D of the calibration factor of no more than 15% for the quantitation runs and 20% for the confirmation runs.

Yes _____ No _____

Comments:

NA.

BLANK ANALYSIS RESULTS BNA, water + soil

The blank analysis were reviewed. The contaminants in the blank are listed below: (*Indicates amounts above CRIL) [Soil - mg/kg]

$$\left[\begin{array}{l} \text{Soil} = \text{g/g} / \text{kg} \\ \text{100 g/m}^3 / \text{ft} \end{array} \right]$$

Remarks: there were other HSL compound present in the blank at levels below the CRL (\approx 0.1 to 0.2 ppb). These are not important and also not confirmed because of their low concentrations.

BLANK ANALYSIS RESULTS - VOA, water & soil

The blank analysis were reviewed. The contaminants in the blank are listed below: (*Indicates amounts above CRDL)

$$\begin{aligned} \text{Soil} &= \mu\text{g}/\text{kg} \\ \text{H}_2\text{O} &= \mu\text{g}/\text{L} \end{aligned}$$

Sample	Date	Compound	Conc.	Samples Associated with this Blank
(MB-1) (VOA/H ₂ O)	4/26	Methylene chloride	2	H 1838 - H 1842
		Acetone	2	
		2-Butanone	1	
(MB-2) (VOA/H ₂ O)	4/27	Methylene chloride	7	
		Acetone	5	
		2-Butanone	1	
		Hexane	8	
		1,1'-Oxybis-ethane	6	H 1838 to H 1842
(MB-3) (VOA/soil)	4/30	Methylene chloride	9	H 1843 - H 1842, H 1849
		Acetone	4	
		2-Butanone	2	
		Toluene	2	
(MB-4) (VOA/soil)	5/1	Methylene chloride	18	
		Acetone	20	
		2-Butanone	2	H 1843 to H 1847, H 1849

Remarks: For MB-1 through MB-3 there were several HSL compounds which were indicated as present in the raw data at a concentration of $\leq 1\text{ ppb}$. These compounds were present below the CRDL and were not listed above for that reason. 2-Butanone was included in MB-2 even though the mass spectral confirmation was not adequate.

DATA COMPLETENESS CHECKLIST

Included; no problems
 Included; problems noted in review
 Not Included

Not Applicable = NA

Case Narrative

Quality Control Summary Package

- Surrogate Recovery Summary (Form II)
- MS/MDS Summary (Form III)
- Reagent Blank Summary (Form IV)
- GC/MS Tuning and Mass Calibration (Form V)

Sample Data Package

- Holding Times (SMO Sample Traffic Reports)
- Organic Analysis Data Sheets (Form I; all four pages for each sample, arranged in increasing SMO number order)
- Reconstructed Ion Chromatogram(s) (RIC)
- GC/EC Chromatograms
- Quantitation Reports
- Mass Spectral Data
- EPA/NIH Mass Spectral Library Search for TIC's

Standards Data Package

- Current List of Laboratory/Instrumental Detection Limits
- Initial Calibration Data (Form VI) for each instrument
- Continuing Calibration Data (Form VII) for each instrument
- Pesticide Evaluation Standards Summary (Form VIII)
- Pesticide/PCB Standards Summary (Form IX)
- Pesticide/PCB Identification (Form X; if any positive results)
- VOA and EPA Standards Reconstruction Ion Chromatograms (RIC)
- VOA and EPA Standards Quantitation Reports
- Pesticide/PCB Standard Chromatograms and Data System Printouts

Raw QC Data Package

- DFTPP and HFB mass spectra and mass listings
- Reagent Blank Data
 - Organic Analysis Data Sheets (Form I)
 - Reconstructed Ion Chromatograms (RIC)
 - Quantitation Reports
 - Mass Spectral Data
 - EPA/NIH Library Search of TIC's
 - GC/EC Chromatograms and Data System Printouts

Matrix Spike and Matrix Spike Duplicate Data

- Organic Analysis Data Sheets (Form I)
- Reconstructed Ion Chromatograms (RIC)
- Quantitation Reports
- Mass Spectral Data
- EPA/NIH Library Search of TIC's
- GC/EC Chromatograms and Data System Printouts

DATA QUALIFIER DEFINITIONS
Region 3

For the purposes of this data review document the following code letters and associated definitions are provided.

- U - The material was analyzed for, but was not detected. The associated numerical value is the estimated sample quantitation limit.
- J - The associated numerical value is an estimated quantity because the amount detected is below the required limits or because quality control criteria were not met.
- UB - Estimated sample quantitation limit increased. Amount found in sample reported. Compound detected at <5 X the amount in blank (<10 X for methylene chloride, acetone, toluene and phthalates).
- W - Detection limit is estimated because quality control criteria were not met.
- JB - The value is an estimated amount detected below required limits and also detected in the blank.
- B - Compound was detected in the blank. Quantity reported is >5 X the amount found in the blank (>10 X for methylene chloride, acetone, toluene, and phthalates).
- R - Quality Control indicates that data is not usable (compound may or may not be present). Resampling and reanalysis is necessary for verification.
- Z - No analytical result.
- N - Presumptive evidence of presence of material (tentative identification).

COMPOUNDS IDENTIFIED

Sample No. H 1838

Hazardous Substances List (HSL) Compounds Detected:

Tentatively Identified Compounds Detected: see page # 3-004

SPECTRUM Matching Quality

- The spectra were examined and found to be of good matching quality.
The spectra were examined and found to be of poor matching quality due to:

REMARKS: Even if the above compounds were not qualified I, as a result of calibration, all of the above were also found in the reagent blanks and would have been qualified up, 3B or B. Because the above compounds were in the blank runs, results should not be considered useful.

COMPOUNDS IDENTIFIED

Sample No. 1939

Hazardous Substances List (HSL) Compounds Detected:

Tentatively Identified Compounds Detected: See page # 3-030

Spectra Matching Quality

- The spectra were examined and found to be of good matching quality.
 The spectra were examined and found to be of poor matching quality due to:

Remarks: see H/128 comments.

COMPOUNDS IDENTIFIED

Sample No. H 1840 water./ Low.

Hazardous Substances List (HSL) Compounds Detected:

Tentatively Identified Compounds Detected: See page # 3-056.

Spectra Matching Quality

- The spectra were examined and found to be of good matching quality.
 The spectra were examined and found to be of poor matching quality due to:

Remarks: See comments H/1938 —

COMPOUNDS IDENTIFIED

Sample No. H 1941 low water

Hazardous Substances List (HSL) Compounds Detected:

Tentatively Identified Compounds Detected: See page # 3-097

Spectra Matching Quality

- The spectra were examined and found to be of good matching quality.
 The spectra were examined and found to be of poor matching quality due to:

Remarks: See comments from sample Miss

COMPOUNDS IDENTIFIED

Sample No. H 1942 Low / water.

Hazardous Substances List (HSL) Compounds Detected:

Tentatively Identified Compounds Detected: See page # 3-11

Spectra Matching Quality

- The spectra were examined and found to be of good matching quality.
 The spectra were examined and found to be of poor matching quality due to:

Remarks: See the comments for sample H1B3B

COMPOUNDS IDENTIFIED

Sample No. H 1943 Low soil

Hazardous Substances List (HSL) Compounds Detected:

Tentatively Identified Compounds Detected: see page 3 - 143

Spectra Matching Quality

- The spectra were examined and found to be of good matching quality. The spectra were examined and found to be of poor matching quality due to:

Remarks: See Comments 4/18/38 -

COMPOUNDS IDENTIFIED

Sample No. H 1344 Low soil

Hazardous Substances List (HSL) Compounds Detected:

Tentatively Identified Compounds Detected: See page # 3-190

Spectra Matching Quality

- The spectra were examined and found to be of good matching quality.
 The spectra were examined and found to be of poor matching quality due to:

Remarks: See comments on sample H1939

COMPOUNDS IDENTIFIED

Sample No. H 1945

Hazardous Substances List (HSL) Compounds Detected:

Tentatively Identified Compounds Detected: see page # 3-226.

Spectra Matching Quality

- The spectra were examined and found to be of good matching quality.
 The spectra were examined and found to be of poor matching quality due to:
Remarks: see comments H1838

Remarks: See comments 4/18/38

COMPOUNDS IDENTIFIED

Sample No. 141846

Hazardous Substances List (HSL) Compounds Detected:

Tentatively Identified Compounds Detected: See page # 3-267

Spectra Matching Quality

- The spectra were examined and found to be of good matching quality.
The spectra were examined and found to be of poor matching quality due to:
See comments on sample H1838

Remarks:

COMPOUNDS IDENTIFIED

Sample No. H 1947

Hazardous Substances List (HSL) Compounds Detected:

Tentatively Identified Compounds Detected: See page # 3-291

Spectra Matching Quality

- The spectra were examined and found to be of good matching quality.
 The spectra were examined and found to be of poor matching quality due to:

COMPOUNDS IDENTIFIED

Sample No. H 1849

Hazardous Substances List (HSL) Compounds Detected:

Tentatively Identified Compounds Detected: See page # 3 - 331.

Spectra Matching Quality

- The spectra were examined and found to be of good matching quality.
 The spectra were examined and found to be of poor matching quality due to:

Remarks: Gounds of Song #1838.

P6
10/18/85

Organics Analysis Data Sheet
(Page 1)

Laboratory Name: CE-EMSI
b Sample ID No: 042685V08 (VOA) 051685C06 (EPA)

Sample Matrix: Water

Data Release Authorized By: Richard J Countess

Case No: 4242
QC Report No: 4242
Contract No: 68-01-6783
Date Sample Received: 24-APR-85

Volatile Compounds

Qualifiers next to compound name are the correct qualifiers.

Concentration: Low
Date Extracted/Prepared: NA
Date Analyzed: 26-APR-85
Conc/Dil Factor: 10.0 pH NA
Percent Moisture: NA
Percent Moisture (Decanted): NA

S
number

71-67-3 Chloromethane
103-9 Bromomethane
75-01-4 Vinyl Chloride J
75-00-3 Chloroethane
109-2 Methylene Chloride J
107-64-1 Acetone J
75-15-0 Carbon Disulfide
103-5-4 1,1-Dichloroethene
103-4-3 1,1-Dichloroethane
156-60-5 Trans-1,2-Dichloroethene
107-66-3 Chloroform
107-06-2 1,2-Dichloroethane
78-93-3 2-Butanone J
71-55-6 1,1,1-Trichloroethane J
103-23-5 Carbon Tetrachloride
103-05-4 Vinyl Acetate R
75-27-4 Bromodichloromethane

S number	ug/l or ug/Kg (Circle One)	CAS Number	ug/l or ug/Kg (Circle One)
71-67-3	200J	79-34-5	1,1,2,2-Tetrachloroethane J
103-9	200	78-67-5	1,2-Dichloropropane J
75-01-4	200L	10061-02-6	Trans-1,3-Dichloropropene
75-00-3	200	79-01-6	Trichloroethene
109-2	1600B	124-48-1	Dibromochloromethane
107-64-1	82B	79-00-5	1,1,2-Trichloroethane
75-15-0	40J	71-43-2	Benzene
103-5-4	300	10061-01-5	cis-1,3-Dichloropropene
103-4-3	200	110-75-8	2-Chloroethylvinylether R
156-60-5	200	75-25-2	Bromoform
107-66-3	100	591-78-6	2-Hexanone J
107-06-2	100	108-10-1	4-Tethyl-2-Pentanone
78-93-3	14 J	127-18-4	Tetrachloroethene
71-55-6	300	108-88-3	Toluene
103-23-5	300	108-90-7	Chlorobenzene
103-05-4	20 R	100-41-4	Ethylbenzene
75-27-4	100	100-42-5	Styrene
			Total Xylenes

Data Reporting Qualifiers

For reporting results to EPA, the following results qualifiers are used.

V Value If the result is a value greater than or equal to the detection limit, report the value

U Compound was analyzed for but not detected.
The number is the minimum attainable detection limit for the sample.

J Indicates an estimated value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicates the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero (e.g. 10J).

C This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides >10 ng/ul in the final extract should be confirmed by GC/MS.

B This flag is used when the analyte is found in the blank as well as a sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action.

S Indicates compound was spiked into the sample for the matrix recovery study. The amounts that were spike into and recovered from the sample are reported on Form III.

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10/11

Organics Analysis Data Sheet
(Page 2)

Qualifiers next
to compound are the
correct qualifiers
except when none is
listed.

Semi-volatile Compounds

Concentration: Low
Date Extracted/Prepared: 26-APR-85
Date Analyzed: 16-MAY-85
Conc/Dil Factor: 1.0

CAS Number	<input type="radio"/> up/l or ug/Kg (Circle One)	CAS Number	<input type="radio"/> up/l or ug/Kg (Circle One)
62-75-9 N-Nitrosodimethylamine	<input checked="" type="radio"/> Correct	201 83-32-9 Acenaphthene	10U
108-95-2 Phenol	<input checked="" type="radio"/> U.J	6U 51-28-5 2,4-Dinitrophenol	7U
2-53-3 Aniline	<input checked="" type="radio"/> U.J	10U 100-02-7 4-Nitrophenol	8U
111-44-4 bis(2-Chloroethyl)Ether	<input checked="" type="radio"/> U.J	10U 132-64-9 Dibenzofuran	8U
95-57-8 2-Chlorophenol	<input type="radio"/>	6U 121-14-2 2,4-Dinitrotoluene	2U
41-73-1 1,3-Dichlorobenzene	<input type="radio"/>	6U 606-20-2 2,6-Dinitrotoluene	2U
66-46-7 1,4-Dichlorobenzene	<input type="radio"/>	6U 84-66-2 Diethylphthalate	10U
100-51-6 Benzyl Alcohol	<input type="radio"/>	20U 7005-72-3 4-Chlorophenyl-phenylether	4U
25-50-1 1,2-Dichlorobenzene	<input type="radio"/>	6U 86-73-7 Fluorene	6U
5-48-7 2-Methylphenol	<input type="radio"/>	6U 100-01-6 4-Nitroaniline	20U
39638-32-9 bis (2-Chloroisopropyl) Ether	<input type="radio"/>	6U 534-52-1 4,6-Dinitro-2-Methylphenol	50U
106-44-5 4-Methylphenol	<input type="radio"/>	30U 86-30-6 N-Nitrosodiphenylamine (1)	2U
21-64-7 N-Nitroso-Di-n-Propylamine	<input type="radio"/>	20U 101-55-3 4-Bromophenylphenylether	10U
57-72-1 Hexachloroethane	<input type="radio"/>	8U 118-74-1 Hexachlorobenzene	8U
98-95-3 Nitrobenzene	<input type="radio"/>	20U 87-86-5 Pentachlorophenol	8U
8-59-1 Isophorone	<input type="radio"/>	10U 25-01-8 Phenanthrene	8U
18-75-5 2-Nitrophenol	<input type="radio"/>	8U 120-12-7 Anthracene	4U
105-67-9 2,4-Dimethylphenol	<input type="radio"/>	4U 84-74-2 Di-n-butylphthalate	4U
65-85-0 Benzoic Acid	<input checked="" type="radio"/> U.J	20U 206-44-0 Fluoranthene	10U
11-91-1 bis (2-Chloroethoxy) Methane	<input type="radio"/>	10U 92-87-5 Benzidine	3U
120-63-2 2,4-Dichlorophenol	<input type="radio"/>	4U 129-00-0 Pyrene	2U
120-82-1 1,2,4-Trichlorobenzene	<input type="radio"/>	10U 85-68-7 Butylbenzylphthalate	10U
11-20-3 Naphthalene	<input type="radio"/>	10U 91-94-1 3,3'-Dichlorobenzidine	2U
106-47-8 4-Chloraniline	<input checked="" type="radio"/> U.J	50U 56-55-3 Benzo(a)anthracene	6U
87-68-3 Hexachlorobutadiene	<input type="radio"/>	10U 117-81-7 bis (2-Ethylhexyl) Phthalate	7
39-50-7 4-Chloro-3-Methylphenol	<input type="radio"/>	8U 218-01-9 Chrysene	6U
31-57-6 2-Methylnaphthalene	<input type="radio"/>	4U 117-84-0 Di-n-octyl Phthalate	6U
77-47-4 Hexachlorocyclopentadiene	<input type="radio"/>	30U 205-99-2 Benzo(b)Fluoranthene	2U
58-06-2 2,4,6-Trichlorophenol	<input type="radio"/>	4U 207-08-9 Benzo(k)Fluoranthene	8U
15-95-4 2,4,5-Trichlorophenol	<input type="radio"/>	4U 50-32-8 Benzo(a)Pyrene	6U
91-58-7 2-Chloronaphthalene	<input type="radio"/>	8U 193-39-5 Indeno(1,2,3-cd)Pyrene	6U
88-74-4 2-Nitroaniline	<input checked="" type="radio"/> R	40U 53-70-3 Dibenzo(a,h)Anthracene	6U
131-11-3 Dimethyl Phthalate	<input type="radio"/>	8U 191-24-2 Benzo(g,h,i)Perylene	6U
206-96-6 Acenaphthylene	<input checked="" type="radio"/> U.J	4U	
99-09-2 3-Nitroaniline	<input checked="" type="radio"/> U.J	40U (1)- Cannot be separated from diphenylaniline	6U

Sample Number

N1838

ANALYSIS NOT REQUIRED

Organics Analysis Data Sheet
(Page 3)

Pesticide/PCBs

Concentration: Low Medium (Circle One)

Date Extracted/Prepared: _____

Date Analyzed: _____

Conc/Dil Factor: _____

CAS Number	ug/l or ug/Kg (Circle One)
319-84-6	Alpha-BHC
319-85-7	Beta-BHC
319-86-8	Delta-BHC
58-89-9	Gamma-BHC (Lindane)
76-44-8	Heptachlor
309-00-2	Aldrin
1024-57-3	Heptachlor Epoxide
959-98-8	Endosulfan I
60-57-1	Dieldrin
72-55-9	4, 4'-DDE
72-20-8	Endrin
33213-65-9	Endosulfan II
72-54-8	4, 4'-DDD
7421-93-4	Endrin Aldehyde
1031-07-8	Endosulfan Sulfate
50-29-3	4, 4'-DDT
72-43-5	Methoxychlor
53494-70-5	Endrin Ketone
57-74-9	Chlordane
8001-35-2	Toxaphene
12674-11-2	Aroclor-1016
11104-28-2	Aroclor-1221
11141-16-5	Aroclor-1232
53469-21-9	Aroclor-1242
12672-29-6	Aroclor-1248
11097-69-1	Aroclor-1254
11096-82-5	Aroclor-1260

V_i = Volume of extract injected (ul)

V_s = Volume of water extracted (ml)

W_s = Weight of sample extracted (g)

V_t = Volume of total extract (ul)

V_s _____ or W_s _____ V_i _____ V_t _____

3 003

Environmental Protection Agency, CLP Sample Management Office,
P O Box 818, Alexandria, Virginia 22313 703/557-2490

Sample Number

H1838

Ab
10/8/84

Organics Analysis Data Sheet
(Page 4)

Tentatively Identified Compounds

No volatile compounds found.

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or ug/kg)
1. --	x,y,z-Trichloro-1-propene	BNA	653	30J
2. --	Unknown	BNA	876	10J
3. --	Unknown polychlorinated compound	BNA	1163	6J
4. --	Unknown amide	BNA	1752	20J
5. --				
6. --				
7. --				
8. --				
9. --				
10. --				
11. --				
12. --				
13. --				
14. --				
15. --				
16. --				
17. --				
18. --				
19. --				
20. --				
21. --				
22. --				
23. --				
24. --				
25. --				
26. --				
27. --				
28. --				
29. --				
30. --				

PL
10/9/85

Organics Analysis Data Sheet
(Page 1)

Laboratory Name: CE-EMSI

Lab Sample ID No: 042685V09 (VOA) 052185C12 (BNA)

Sample Matrix: Water

Data Release Authorized By:

Richard J Courtney

Case No: 4242

QC Report No: 4242

Contract No: 68-01-6783

Date Sample Received: 24-APR-85

Volatile Compounds

Concentration: Low
Date Extracted/Prepared: NA
Date Analyzed: 26-APR-85
Conc/Dil Factor: NA pH NA
Percent Moisture: NA
Percent Moisture (Decanted): NA

IS
Number

u/l or ug/kg
(Circle One)

CAS
Number

u/l or ug/kg
(Circle One)

4-87-3 Chloromethane
4-83-9 Bromomethane
75-01-4 Vinyl Chloride WJ
75-00-3 Chloroethane
5-09-2 Methylene Chloride J
67-64-1 Acetone J
75-15-0 Carbon Disulfide
5-35-4 1,1-Dichloroethene
75-34-3 1,1-Dichloroethane
156-60-5 Trans-1,2-Dichloroethene
7-66-3 Chloroform
07-06-2 1,2-Dichloroethane
78-93-3 2-Butanone J
71-55-6 1,1,1-Trichloroethane WJ
6-23-5 Carbon Tetrachloride
108-05-4 Vinyl Acetate R
75-27-4 Bromodichloromethane

20U	79-34-5	1,1,2,2-Tetrachloroethane WJ	WJ
2U	78-87-5	1,2-Dichloropropane WJ	WJ
20U	10061-02-6	Trans-1,3-Dichloropropene	WJ
2U	79-01-6	Trichloroethene	WJ
300B	124-48-1	Dibromochloromethane	J
38B	79-00-5	1,1,2-Trichloroethane	J
4U	71-43-2	Benzene	WJ
3U	10061-01-5	cis-1,3-Dichloropropene	WJ
2U	110-75-8	2-Chloroethylvinylether R	WJ
2U	75-25-2	Bromoform	J
1U	591-78-6	2-Hexanone WJ	J
1U	108-10-1	4-Methyl-2-Pentanone	WJ
420B	127-18-4	Tetrachloroethene	WJ
3U	108-88-3	Toluene	WJ
3U	108-90-7	Chlorobenzene	57
2U	100-41-4	Ethylbenzene	WJ
1U	100-42-5	Styrene	WJ
Total Xylenes			

Data Reporting Qualifiers

For reporting results to EPA, the following results qualifiers are used.

V Value If the result is a value greater than or equal to the detection limit, report the value

W Compound was analyzed for but not detected.
The number is the minimum attainable detection limit for the sample.

J Indicates an estimated value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicates the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero (e.g. 10J).

R This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides $>10 \text{ ng/uL}$ in the final extract should be confirmed by GC/MS.

B This flag is used when the analyte is found in the blank as well as a sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action.

S Indicates compound was spiked into the sample for the matrix recovery study. The amounts that were spiked into and recovered from the sample are reported on Form III.

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10/9/85

Organics Analysis Data Sheet
(Page 2)

Semi-volatile Compounds

Concentration: Low
Date Extracted/Prepared: 17-MAY-85
Date Analyzed: 21-MAY-85 ←
Conc/Dil Factor: 1.0

Holding time exceeded for extraction
re-analysis

CAS Number	ug/l or ug/kg (Circle One)	CAS Number	ug/l or ug/kg (Circle One)
62-75-9	N-Nitrosodimethylamine	20U	83-32-9
108-95-2	Phenol - UJ	6U	51-28-5
62-53-3	Aniline - UJ	10U	100-02-7
111-44-4	bis(2-Chloroethyl)Ether-UJ	10U	132-64-9
95-57-8	2-Chlorophenol	6U	121-14-2
541-73-1	1,3-Dichlorobenzene	6U	606-20-2
106-46-7	1,4-Dichlorobenzene	6U	84-66-2
100-51-6	Benzyl Alcohol	20U	7005-72-3
95-50-1	1,2-Dichlorobenzene	6U	86-73-7
95-48-7	2-Methylphenol	6U	100-01-6
39638-32-9	bis (2-Chloroisopropyl) Ether	6U	534-52-1
106-44-5	4-Methylphenol	30U	86-30-6
621-64-7	N-Nitroso-Di-n-Propylamine	20U	101-55-3
67-72-1	Hexachloroethane	8U	118-74-1
98-95-3	Nitrobenzene	20U	87-86-5
78-59-1	Isophorone	10U	85-01-8
88-75-5	2-Nitrophenol	8U	120-12-7
105-67-9	2,4-Dimethylphenol	4U	84-74-2
65-85-0	Benzoic Acid - UJ	20U	206-44-0
111-91-1	bis (2-Chloroethoxy) Methane	10U	92-87-5
120-83-2	2,4-Dichlorophenol	4U	129-00-0
120-82-1	1,2,4-Trichlorobenzene	10U	85-68-7
91-20-3	Naphthalene	10U	91-94-1
106-47-8	4-Chloraniline-UJ	50U	56-55-3
87-68-3	Hexachlorobutadiene	10U	117-81-7
59-50-7	4-Chloro-3-Methylphenol	8U	218-01-9
91-57-6	2-Methylnaphthalene	4U	117-84-0
77-47-4	Hexachlorocyclopentadiene	30U	205-99-2
88-06-2	2,4,6-Trichlorophenol	4U	207-08-9
95-95-4	2,4,5-Trichlorophenol	4U	50-32-8
91-58-7	2-Chloronaphthalene	8U	193-39-5
38-74-4	2-Nitroaniline - R	40U	53-70-3
131-11-3	Dimethyl Phthalate	8U	191-24-2
108-96-8	Acenaphthylene - UJ	4U	
109-09-2	3-Nitroaniline. UJ	40U	

(1) - Cannot be separated from diphenylamine

Sample Number

H1839

PL
10/27/84

Organics Analysis Data Sheet
(Page 3)

ANALYSIS NOT REQUIRED

Pesticide/PCBs

Concentration: Low Medium (Circle One)

Date Extracted/Prepared: _____

Date Analyzed: _____

Conc/Dil Factor: _____

CAS Number		ug/l or ug/Kg (Circle One)
319-84-6	Alpha-BHC	
319-85-7	Beta-BHC	
319-86-8	Delta-BHC	
58-89-9	Gamma-BHC (Lindane)	
76-44-8	Heptachlor	
309-00-2	Aldrin	
1024-67-3	Heptachlor Epoxide	
959-98-8	Endosulfan I	
60-57-1	Dieldrin	
72-55-9	4, 4'-DDE	
72-20-8	Endrin	
23213-65-9	Endosulfan II	
72-54-8	4, 4'-DDD	
7421-93-4	Endrin Aldehyde	
1031-07-8	Endosulfan Sulfate	
50-29-3	4, 4'-DDT	
72-43-5	Methoxychlor	
53494-70-5	Endrin Ketone	
57-74-9	Chlordane	
8001-35-2	Toxaphene	
12674-11-2	Aroclor-1016	
11104-28-2	Aroclor-1221	
11141-16-5	Aroclor-1232	
53469-21-9	Aroclor-1242	
12672-29-6	Aroclor-1248	
11097-69-1	Aroclor-1254	
11096-82-9	Aroclor-1260	

V_1 = Volume of extract injected (ul)

V_s = Volume of water extracted (ml)

W_s = Weight of sample extracted (g)

V_t = Volume of total extract (ul)

V_s _____ or W_s _____ V_1 _____ V_t _____

Environmental Protection Agency CLP Sample Management Office
P O Box 818, Alexandria, Virginia 22313 703/557-2490

Sample Number
H1839

Pb
10/18/85

Organics Analysis Data Sheet
(Page 4)

Tentatively Identified Compounds

No volatile compounds found.

CAS Number	Compound Name	Fraction	RT on Scan Number	Estimated Concentration (ug/l or ug/kg)
1. --	x,y,z-Trichloro-1-propene	BNA	637	10JB
2. --	Unknown	BNA	852	4J
3.				
4.				
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3 030
4/84

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Box 616. Alexandria, Virginia 22313 703/557-2490

Sample Number
H1840

Notice. It is due to the
quality of the document
being filed

10/16/85
RJ

Organics Analysis Data Sheet
(Page 1)

Laboratory Name: CE-EMSI
Sample ID No: 042685V13 (VOA) 051685C11 (EPA)
Sample Matrix: Water
Data Release Authorized By: Richard J Counters

Case No: 4242
QC Report No: 4242
Contract No: 68-01-6783
Date Sample Received: 24-APR-85

Volatile Compounds

Concentration: Low
Date Extracted/Prepared: NA
Date Analyzed: 26-APR-85
Conc/Dil Factor: NA pH NA
Percent Moisture: NA
Percent Moisture (Decanted): NA

I S number	ug/l or ug/Kg (Circle One)	CAS Number	ug/l or ug/Kg (Circle One)
-87-3	Chloromethane	20U	79-34-5 1,1,2,2-Tetrachloroethane - U J
-83-9	Bromomethane	2U	78-87-5 1,2-Dichloropropane -- U J
75-01-4	Vinyl Chloride - U J	20U	10061-02-6 Trans-1,3-Dichloropropene
75-00-3	Chloroethane	2U	79-01-6 Trichloroethene
-09-2	Methylene Chloride - J	12B	124-48-1 Dibromochloromethane
67-64-1	Acetone - J	140B	79-00-5 1,1,2-Trichloroethane
75-15-0	Carbon Disulfide	4U	71-43-2 Benzene
-35-4	1,1-Dichloroethene	3U	10061-01-5 cis-1,3-Dichloropropene
72-34-3	1,1-Dichloroethane	2U	110-75-8 2-Chloroethylvinylether - R
156-60-5	Trans-1,2-Dichloroethene	2U	75-25-2 Bromoform
-66-3	Chloroform	1U	591-78-6 2-Hexanone - U J
77-06-2	1,2-Dichloroethene	1U	108-10-1 4-Methyl-2-Pentanone
78-93-3	2-Butanone - J	200B	127-18-4 Tetrachloroethene
77-55-6	1,1,1-Trichloroethane - U J	3U	108-88-3 Toluene
-23-5	Carbon Tetrachloride	3U	108-90-7 Chlorobenzene
108-05-4	Vinyl Acetate - R	2U	100-41-4 Ethylbenzene
75-27-4	Bromodichloromethane	1U	100-42-5 Styrene
			Total Xylenes

Data Reporting Qualifiers

For reporting results to EPA, the following results qualifiers are used.

V Value If the result is a value greater than or equal to the detection limit, report the value

C This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides >10 ng/ul in the final extract should be confirmed by GC/MS.

B This flag is used when the analyte is found in the blank as well as a sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action.

S Indicates compound was spiked into the sample for the matrix recovery study. The amounts that were spiked into and recovered from the sample are reported on Form III.

J Indicates an estimated value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicates the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero (e.g. 10U).

SITE / SITE ASSESSMENT
SASOOG

26
2/84

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20 Box 818. Alexandria, Virginia 22313 703/557-2480

notice, it is due to the quality of the document being filmed

SITE / SITE ASSESSMENT

SAS006

E 0 4 7

Sample Number
H1840

Organics Analysis Data Sheet
(Page 2)

Semi-volatile Compounds

Concentration: Low
Date Extracted/Prepared: 26-APR-85
Date Analyzed: 16-MAY-85
Conc/Dil Factor: 1.0

CAS Number	ug/l or ug/Kg (Circle One)	CAS Number	ug/l or ug/Kg (Circle One)
62-75-9	N-Nitrosodimethylamine	20U	83-32-9
108-95-2	Phenol - uS	6U	51-28-5
2-53-3	Aniline - uJ	10U	100-02-7
11-44-4	bis(2-Chloroethyl)Ether - uJ	10U	132-64-9
95-57-8	2-Chlorophenol	6U	121-14-2
41-73-1	1,3-Dichlorobenzene	6U	606-20-2
36-46-7	1,4-Dichlorobenzene	6U	84-66-2
100-51-6	Benzyl Alcohol	20U	7005-72-3
95-50-1	1,2-Dichlorobenzene	6U	86-73-7
5-48-7	2-Methylphenol	6U	100-01-6
39638-32-9	bis (2-Chloroisopropyl) Ether	6U	534-52-1
106-44-5	4-Methylphenol	30U	86-30-6
21-64-7	N-Nitroso-Di-n-Propylamine	20U	101-55-3
7-72-1	Hexachloroethane	8U	118-74-1
98-95-3	Nitrobenzene	20U	87-86-5
78-59-1	Isophorone	10U	85-01-8
8-75-5	2-Nitrophenol	8U	120-12-7
105-67-9	2,4-Dimethylphenol	4U	84-74-2
65-85-0	Benzoic Acid - uJ	20U	206-44-0
11-91-1	bis (2-Chloroethoxy) Methane	10U	92-87-5
220-83-2	2,4-Dichlorophenol	4U	129-00-0
120-82-1	1,2,4-Trichlorobenzene	10U	85-68-7
1-20-3	Naphthalene	10U	91-24-1
06-47-8	4-Chloroaniline - uJ	50U	56-55-3
87-68-3	Hexachlorobutadiene	10U	117-81-7
59-50-7	4-Chloro-3-Methylphenol	8U	218-01-9
1-57-6	2-Methylnaphthalene	4U	117-84-0
77-47-4	Hexachlorocyclopentadiene	30U	205-99-2
88-06-2	2,4,6-Trichlorophenol	4U	207-08-9
5-95-4	2,4,5-Trichlorophenol	4U	50-32-8
1-58-7	2-Chloronaphthalene	8U	193-39-5
88-74-4	2-Nitroaniline - R	40U	53-70-3
31-11-3	Dimethyl Phthalate	8U	191-24-2
08-96-8	Acenaphthylene ~ uS	4U	
99-09-2	3-Nitroaniline ~ uJ	40U	

(1) Cannot be separated from diphenylamine

Environmental Protection Agency, CLP Sample Management Office,
P. O. Box 810, Alexandria, Virginia 22313 703/557-2490

Sample Number

H1840

RE
AF/BS

ANALYSIS NOT REQUIRED

Organics Analysis Data Sheet
(Page 3)

Pesticide/PCBs

Concentration: Low Medium (Circle One)

Date Extracted/Prepared: _____

Date Analyzed: _____

Conc/Dil Factor: _____

CAS Number	ug/l or ug/Kg (Circle One)
------------	----------------------------

319-84-6	Alpha-BHC
319-85-7	Beta-BHC
319-86-8	Delta-BHC
58-89-9	Gamma-BHC (Lindane)
76-44-8	Heptachlor
309-00-2	Aldrin
1024-57-3	Heptachlor Epoxide
959-98-8	Endosulfan I
60-57-1	Dieldrin
72-55-9	4, 4'-DDE
72-20-8	Endrin
33213-65-9	Endosulfan II
72-54-8	4, 4'-DDD
7421-93-4	Endrin Aldehyde
1031-07-8	Endosulfan Sulfate
50-29-3	4, 4'-DDT
72-43-5	Methoxychlor
53494-70-5	Endrin Ketone
57-74-9	Chlordane
8001-35-2	Toxaphene
12674-11-2	Aroclor-1016
11104-28-2	Aroclor-1221
11141-16-5	Aroclor-1232
53469-21-9	Aroclor-1242
12672-29-6	Aroclor-1248
11097-69-1	Aroclor-1254
11096-82-5	Aroclor-1260

V_i = Volume of extract injected (ul)

V_s = Volume of water extracted (ml)

W_s = Weight of sample extracted (g)

V_t = Volume of total extract (ul)

V_s _____ or W_s _____ V_t _____ V_i _____

3 055

Environmental Protection Agency, CLP Sample Management Office,
P O Box 818, Alexandria, Virginia 22313 703/557-2480

Sample Number
H1840

PP
0/7/85

Organics Analysis Data Sheet
(Page 4)

Tentatively Identified Compounds

No volatile compounds found.

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or ug/kg)
1. --	X, Y, Z-Trichloro-1-propene	BNA	656	8J
2. --	Unknown saturated aliphatic hydrocarbon	BNA	1327	5J
3. --	Unknown	BNA	1373	4J
4. --	Unknown	BNA	1468	8J
5. --	Unknown amide	BNA	1747	15J
6. --				
7. --				
8. --				
9. --				
10. --				
11. --				
12. --				
13. --				
14. --				
15. --				
16. --				
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23. --				
24. --				
25. --				
26. --				
27. --				
28. --				
29. --				
30. --				

Environmental Protection Agency, CLP Sample Management Office,
P.O. Box 816, Alexandria, Virginia 22313 703/557-2490

Sample Number
H1841

26
10/11/85

Organics Analysis Data Sheet
(Page 1)

Laboratory Name: CE-EMSI

Sample ID No: 042783V11 (VOA) 051685C12 (BNA)

Sample Matrix: Water

Data Release Authorized By:

Richard Lourdes

Case No: 4242

QC Report No: 4242

Contract No: 68-01-6783

Date Sample Received: 24-APR-85

Volatile Compounds

Concentration: Low
Date Extracted/Prepared: NA
Date Analyzed: 27-APR-85
Conc/Dil Factor: NA pH NA
Percent Moisture: NA
Percent Moisture (Decanted): NA

S number	ug/l or ug/Kg (Circle One)	CAS Number	ug/l or ug/Kg (Circle One)
-97-3	Chloromethane	201	79-34-5
-33-9	Bromomethane	21	78-87-5
75-01-4	Vinyl Chloride u J	201	10061-02-6
75-00-3	Chloroethane	21	79-01-6
109-2	Methylene Chloride J	68	124-48-1
67-64-1	Acetone u J	111	79-00-5
75-15-0	Carbon Disulfide	41	71-43-2
1035-4	1,1-Dichloroethene	31	10061-01-5
1034-3	1,1-Dichloroethane	21	110-75-8
156-60-5	Trans-1,2-Dichloroethene	21	75-25-2
106-63-3	Chloroform	111	591-78-6
107-06-2	1,2-Dichloroethane	111	108-10-1
78-93-3	2-Butanone - R	111	127-18-4
71-55-6	1,1,1-Trichloroethane u J	31	108-88-3
523-5	Carbon Tetrachloride	31	108-90-7
108-05-4	Vinyl Acetate R	21	100-41-4
75-27-4	Bromodichloromethane	111	100-42-5
			Total Xylenes

Data Reporting Qualifiers

For reporting results to EPA, the following results qualifiers are used.

V Value If the result is a value greater than or equal to the detection limit, report the value

U Compound was analyzed for but not detected. The number is the minimum attainable detection limit for the sample.

J Indicates an estimated value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicates the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero (e.g. 10J).

C This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides $\geq 10 \text{ ng}/\text{ul}$ in the final extract should be confirmed by GC/MS.

B This flag is used when the analyte is found in the blank as well as a sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action.

S Indicates compound was spiked into the sample for the matrix recovery study. The amounts that were spiked into and recovered from the sample are reported on Form III.

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Sample Number
H1841

7/3
10/1984
being final

Organics Analysis Data Sheet
(Page 2)

Semi-volatile Compounds

Concentration: Low
Date Extracted/Prepared: 26-APR-85
Date Analyzed: 16-MAY-85
Conc/Dil Factor: 1.0

CAS Number	(<input checked="" type="radio"/> ug/l or ug/kg (Circle One))	CAS Number	(<input checked="" type="radio"/> ug/l or ug/kg (Circle One))
62-75-9	N-Nitrosodimethylamine	20U	63-32-9
68-95-2	Phenol - <input checked="" type="checkbox"/> R	6U	51-28-5
2-53-3	Aniline - <input checked="" type="checkbox"/> J	10U	100-02-7
111-44-4	bis(2-Chloroethyl)Ether <input checked="" type="checkbox"/> J	10U	132-64-9
75-57-6	2-Chlorophenol	6U	121-14-2
41-73-1	1,3-Dichlorobenzene	6U	606-20-2
106-46-7	1,4-Dichlorobenzene	6U	84-66-2
100-51-6	Benzyl Alcohol	20U	7005-72-3
5-50-1	1,2-Dichlorobenzene	6U	86-73-7
5-48-7	2-Methylphenol	6U	100-01-6
39638-32-9	bis (2-Chloroisopropyl) Ether	6U	534-52-1
66-44-5	4-Methylphenol	30U	86-30-6
21-64-7	N-Nitroso-Di-n-Propylamine	20U	101-55-3
67-72-1	Hexachloroethane	8U	118-74-1
98-95-3	Nitrobenzene	20U	67-86-5
8-59-1	Isophorone	10U	85-01-8
88-75-5	2-Nitrophenol	8U	120-12-7
105-67-9	2,4-Dimethylphenol	4U	84-74-2
5-85-0	Benzoic Acid <input checked="" type="checkbox"/> J	20U	206-44-0
111-91-1	bis (2-Chloroethyl) Methane	10U	92-87-5
120-83-2	2,4-Dichlorophenol	4U	129-00-0
120-82-1	1,2,4-Trichlorobenzene	10U	85-68-7
71-20-3	Naphthalene	10U	91-94-1
106-47-8	4-Chloroaniline <input checked="" type="checkbox"/> J	50U	56-55-3
87-68-3	Hexachlorobutadiene	10U	117-81-7
59-50-7	4-Chloro-3-Methylphenol	8U	218-01-9
91-57-6	2-Methylnaphthalene	4U	117-84-0
77-47-4	Hexachlorocyclopentadiene	30U	205-99-2
38-06-2	2,4,6-Trichlorophenol	4U	207-08-9
35-95-4	2,4,5-Trichlorophenol	4U	50-32-8
91-58-7	2-Chloronaphthalene	8U	193-39-5
98-74-4	2-Nitroaniline - <input checked="" type="checkbox"/> R	40U	53-70-3
131-11-3	Dimethyl Phthalate	8U	191-24-2
208-96-8	Acenaphthylene <input checked="" type="checkbox"/> J	4U	Benzo(g,h,i)Perylene
99-09-2	3-Nitroaniline <input checked="" type="checkbox"/> J	40U	(1)- Cannot be separated from diphenylaniline

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26
10/16/97

Sample Number

H1841

ANALYSIS NOT REQUIRED

Organics Analysis Data Sheet
(Page 3)

Pesticide/PCBs

Concentration: Low Medium (Circle One)

Date Extracted/Prepared: _____

Date Analyzed: _____

Conc/Dil Factor: _____

CAS Number	ug/l or ug/Kg (Circle One)
------------	----------------------------

319-84-6	Alpha-BHC
319-85-7	Beta-BHC
319-86-8	Delta-BHC
58-89-9	Gamma-BHC (Lindane)
76-44-8	Heptachlor
309-00-2	Aldrin
1024-57-3	Heptachlor Epoxide
959-98-8	Endosulfan I
60-57-1	Dieldrin
72-55-9	4, 4'-DDE
72-20-8	Endrin
33213-65-9	Endosulfan II
72-54-8	4, 4'-DDD
7421-93-4	Endrin Aldehyde
1031-07-8	Endosulfan Sulfate
50-29-3	4, 4'-DDT
72-43-5	Methoxychlor
53494-70-5	Endrin Ketone
57-74-9	Chlordane
8001-35-2	Toxaphene
12674-11-2	Aroclor-1016
11104-28-2	Aroclor-1221
11141-16-5	Aroclor-1232
53469-21-9	Aroclor-1242
12672-29-6	Aroclor-1248
11097-69-1	Aroclor-1254
11096-82-5	Aroclor-1260

V_i = Volume of extract injected (ul)

V_s = Volume of water extracted (ml)

W_s = Weight of sample extracted (g)

V_t = Volume of total extract (ul)

V_s _____ or W_s _____ V_t _____ V_i _____

Sample Number
H1841

7/2/84

Organics Analysis Data Sheet
(Page 4)

Tentatively Identified Compounds

No volatile compounds found.

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or ug/kg)
1. --	x,y,z-Trichloro-1-propene	BNA	651	6J
2. --	Unknown saturated aliphatic hydrocarbon	BNA	1326	7J
3. --	Unknown	BNA	1467	7J
4. --	Unknown amide	BNA	1745	10J
5.				
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Environmental Protection Agency, CLP Sample Management Office,
P O Box 818, Alexandria, Virginia 22313 703/557-2490

Sample Number
H1842

3/25
108

Organics Analysis Data Sheet
(Page 1)

Laboratory Name: GE-DSI

Lab Sample ID No: 042785V07 (VOA) 052185C03 (EPA)

Sample Matrix: Water

Data Release Authorized By: Richard J. ConTess

Case No: 4242

QC Report No: 4242

Contract No: 68-01-6783

Date Sample Received: 24-APR-85

Volatile Compounds

Concentration: Low
Date Extracted/Prepared: NA
Date Analyzed: 27-APR-85
Conc/Dil Factor: 2. pH NA
Percent Moisture: NA
Percent Moisture (Decanted): NA

CAS
Number

ug/l or ug/kg
(Circle One)

CAS
Number

ug/l or ug/kg
(Circle One)

-67-3	Chloromethane	500	79-34-5	1,1,2,2-Tetrachloroethane <input checked="" type="checkbox"/>	6
74-83-9	Bromomethane	40	78-87-5	1,2-Dichloropropane <input checked="" type="checkbox"/>	2
75-01-4	Vinyl Chloride <input checked="" type="checkbox"/>	500	10061-02-6	Trans-1,3-Dichloropropene	2
500-3	Chloroethane	40	79-01-6	Trichloroethene	4
75-09-2	Methylene Chloride <input checked="" type="checkbox"/>	22B	124-48-1	Dibromoform	6
67-64-1	Acetone <input checked="" type="checkbox"/>	1400B	79-00-5	1,1,2-Trichloroethane	6
5-15-0	Carbon Disulfide	80	71-43-2	Benzene	2
5-35-4	1,1-Dichloroethene	60	10061-01-5	cis-1,3-Dichloropropene	2
75-34-3	1,1-Dichloroethane	40	110-75-8	2-Chloroethylvinylether <input checked="" type="checkbox"/>	100
56-60-5	Trans-1,2-Dichloroethene	40	75-25-2	Bromoform	6
7-66-3	Chloroform	20	591-78-6	2-Hexanone <input checked="" type="checkbox"/>	2
107-06-2	1,2-Dichloroethene	20	108-10-1	4-Methyl-2-Pentanone	2
78-93-3	2-Butanone <input checked="" type="checkbox"/>	22	127-18-4	Tetrachloroethene	2
1-55-6	1,1,1-Trichloroethane <input checked="" type="checkbox"/>	60	108-88-3	Toluene	6
56-23-5	Carbon Tetrachloride	60	108-90-7	Chlorobenzene	6
108-05-4	Vinyl Acetate <input checked="" type="checkbox"/>	40	100-41-4	Ethylbenzene	4
5-27-4	Bromodichloromethane	20	100-42-5	Styrene	4
				Total Xylenes	4

Data Reporting Qualifiers

For reporting results to EPA, the following results qualifiers are used.

A Value If the result is a value greater than or equal to the detection limit, report the value

B Compound was analyzed for but not detected.

The number is the minimum attainable detection limit for the sample.

J Indicates an estimated value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicates the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero (e.g. 10).

C This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides $\geq 10 \text{ ng}/\text{ul}$ in the final extract should be confirmed by GC/MS.

B This flag is used when the analyte is found in the blank as well as a sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action.

S Indicates compound was spiked into the sample for the matrix recovery study. The amounts that were spiked into and recovered from the sample are reported on Form III.

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6/1/85
being run

Organics Analysis Data Sheet
(Page 2)

Semi-volatile Compounds

Concentration: Low
Date Extracted/Prepared: 17-MAY-85
Date Analyzed: 21-MAY-85
Conc/Dil Factor: 1.5

holding time exceeded for extraction.

CAS Number	(ug/l) or ug/Kg (Circle One)	CAS Number	(ug/l) or ug/Kg (Circle One)
62-75-9	N-Nitrosodimethylamine	30U	83-32-9
1-3-95-2	Phenol - u J	9B	51-28-5
1-53-3	Aniline - u J	20U	100-02-7
111-44-4	bis(2-Chloroethyl)Ether u J	20U	132-64-9
1-57-8	2-Chlorophenol	9U	121-14-2
1-73-1	1,3-Dichlorobenzene	9U	606-20-2
106-46-7	1,4-Dichlorobenzene	9U	84-66-2
100-51-6	Benzyl Alcohol	30U	7005-72-3
1-50-1	1,2-Dichlorobenzene	9U	86-73-7
1-48-7	2-Methylphenol	9U	100-01-6
39638-32-9	bis (2-Chloroisopropyl) Ether	9U	534-52-1
1-44-5	4-Methylphenol	40U	86-30-6
1-64-7	N-Nitroso-Di-n-Propylamine	40U	101-55-3
67-72-1	Hexachloroethane	10U	118-74-1
10-95-3	Nitrobenzene	40U	67-86-5
-59-1	Isophorone	20U	85-01-8
88-75-5	2-Nitrophenol	10U	120-12-7
105-67-9	2,4-Dinethylphenol	6U	84-74-2
-85-0	Benzoic Acid u J	30U	206-44-0
111-91-1	bis (2-Chloroethoxy) Methane	20U	92-87-5
120-63-2	2,4-Dichlorophenol	6U	129-00-0
10-82-1	1,2,4-Trichlorobenzene	20U	85-68-7
-20-3	Naphthalene	20U	91-94-1
106-47-8	4-Chloroaniline u J	70U	56-55-3
97-68-3	Hexachlorobutadiene	20U	117-81-7
1-50-7	4-Chloro-3-Methylphenol	10U	218-01-9
91-57-6	2-Methylnaphthalene	6U	117-84-0
77-47-4	Hexachlorocyclopentadiene	40U	205-99-2
1-06-2	2,4,6-Trichlorophenol	6U	207-08-9
1-95-4	2,4,5-Trichlorophenol	6U	50-32-8
91-58-7	2-Chloronaphthalene	10U	193-39-5
3-74-4	2-Nitroaniline - R	60U	53-70-3
31-11-3	Dimethyl Phthalate	10U	191-24-2
208-96-8	Acenaphthylene - u J	6U	(1) - Cannot be separated from diphenylamine
99-09-2	3-Nitroaniline - u J	60U	

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Sample Number

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H1842

ANALYSIS NOT REQUIRED

Organics Analysis Data Sheet
(Page 3)

Pesticide/PCBs

Concentration: Low Medium (Circle One)

Date Extracted/Prepared: _____

Date Analyzed: _____

Conc/Dil Factor: _____

CAS Number		ug/l or ug/Kg (Circle One)
319-84-6	Alpha-BHC	
319-85-7	Beta-BHC	
319-86-8	Delta-BHC	
58-89-9	Gamma-BHC (Lindane)	
76-44-8	Heptachlor	
309-00-2	Aldrin	
1024-57-3	Heptachlor Epoxide	
959-98-8	Endosulfan I	
60-57-1	Dieldrin	
72-55-9	4, 4'-DDE	
72-20-8	Endrin	
33213-65-9	Endosulfan II	
72-54-8	4, 4'-DDD	
7421-93-4	Endrin Aldehyde	
1031-07-8	Endosulfan Sulfate	
50-29-3	4, 4'-DDT	
72-43-5	Methoxychlor	
53494-70-5	Endrin Ketone	
57-74-9	Chlordane	
8001-35-2	Toxaphene	
12674-11-2	Aroclor-1016	
11104-28-2	Aroclor-1221	
11141-16-5	Aroclor-1232	
53489-21-9	Aroclor-1242	
12672-29-6	Aroclor-1248	
11097-69-1	Aroclor-1254	
11096-82-5	Aroclor-1260	

V_i = Volume of extract injected (ul)

V_s = Volume of water extracted (ml)

W_s = Weight of sample extracted (g)

V_t = Volume of total extract (ul)

V_s _____ or W_s _____ V_t _____ V_i _____

Sample Number

H1842

74
6/2/84

Organics Analysis Data Sheet
(Page 4)

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or ug/kg)
1. 109-99-9	Furan, tetrahydro	VOA	93	20J
2. 110-54-3	Hexane	VOA	239	10JB
3. --	x,y,z-Trichloro-1-propene	BNA	637	90JB
4. --	Unknown	BNA	851	40J
5. 1541-20-4	E1-2-cyclohexen-1-yl	BNA	1028	7J
6. --	Unknown polyhalogenated compound	BNA	1132	10JB
7.				
8.	-			
9.				
10.				
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12.				
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P.O. Box 818, Alexandria, Virginia 22313 703/557-2490

Sample Number
H1843

2
alpha

Organics Analysis Data Sheet
(Page 1)

Laboratory Name: CE-EMSI
Lab Sample ID No: 04308SV12 (VOA) 052285C14 (BNA)
Sample Matrix: Soil
Lab Release Authorized By: Richard J Courtney

Case No: 4242
QC Report No: 4242
Contract No: 68-01-6783
Date Sample Received: 24-APR-85

Volatile Compounds

Concentration: Low
Date Extracted/Prepared: NA
Date Analyzed: 30-APR-85
Conc/Dil Factor: NA pH 7.45
Percent Moisture: 34.58
Percent Moisture (Decanted): NA

Number	ug/l or ug/kg (Circle one)	CAS Number	ug/l or ug/kg (Circle one)
-87-3	Chloromethane	40	79-34-5
74-83-9	Bromomethane	30	78-87-5
75-01-4	Vinyl Chloride ^{WJ}	400	10061-02-6
-00-3	Chloroethane	30	79-01-6
-09-2	Methylene Chloride ^J	128	124-48-1
67-64-1	Acetone ^J	223	79-00-5
-15-0	Carbon Disulfide	60	71-43-2
-35-4	1,1-Dichloroethene	50	10061-01-5
75-34-3	1,1-Dichloroethane	30	110-75-8
156-60-5	Trans-1,2-Dichloroethene	30	75-25-2
-66-3	Chloroform	20	591-78-6
707-06-2	1,2-Dichloroethane	20	108-10-1
78-93-3	2-Butanone ^J	68	127-18-4
-55-6	1,1,1-Trichloroethane ^{WJ}	30	108-88-3
-23-5	Carbon Tetrachloride	30	108-90-7
108-05-4	Vinyl Acetate ^R	30	100-41-4
75-27-4	Bromodichloromethane	20	100-42-5
			Total Xylenes

Data Reporting Qualifiers

For reporting results to EPA, the following results qualifiers are used.

Value If the result is a value greater than or equal to the detection limit, report the value

C This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides >10 ng/ μ l in the final extract should be confirmed by GC/MS.

U Compound was analyzed for but not detected. The number is the minimum attainable detection limit for the sample.

B This flag is used when the analyte is found in the blank as well as a sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action.

J Indicates an estimated value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicates the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero (e.g. 10J).

S Indicates compound was spiked into the sample for the matrix recovery study. The amounts that were spiked into and recovered from the sample are reported on Form III.

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Sample Number
N1843

7/2
reptile

being filmed

Organics Analysis Data Sheet
(Page 2)

Semi-volatile Compounds

Concentration: Low
Date Extracted/Prepared: 30-APR-85
Date Analyzed: 22-MAY-85
Conc/Dil Factor: 2.0

CAS Number	ug/l or ug/Kg (Circle One)	CAS Number	ug/l or ug/kg (Circle One)
62-75-9 N-Nitrosodimethylamine	1000U	63-32-9 Acenaphthene	600U
8-95-2 Phenol J S	400U	51-28-5 2,4-Dinitrophenol - U J	4000U
1-53-3 Aniline - U J	800U	100-02-7 4-Nitrophenol - U J	5000U
111-44-4 bis(2-Chloroethyl)Ether U J	800U	132-64-9 Dibenzofuran	500U
52-57-8 2-Chlorophenol	400U	121-14-2 2,4-Dinitrotoluene	2000U
1-73-1 1,3-Dichlorobenzene	400U	606-20-2 2,6-Dinitrotoluene	1000U
106-46-7 1,4-Dichlorobenzene	400U	84-66-2 Diethylphthalate	600U
100-51-6 Benzyl Alcohol	1000U	7005-72-3 4-Chlorophenyl-phenylether	300U
50-1 1,2-Dichlorobenzene	400U	86-73-7 Fluorene	400U
53-48-7 2-Methylphenol	400U	100-01-6 4-Nitroaniline - U J	2000U
39638-32-9 bis (2-Chloroisopropyl) Ether	400U	534-52-1 4,6-Dinitro-2-Methylphenol	3000U
6-44-5 4-Methylphenol	2000U	86-30-6 N-Nitrosodiphenylamine (1)	1000U
1-64-7 N-Nitroso-Di-n-Propylamine	2000U	101-55-3 4-Bromoethylphenylether - U J	900U
67-72-1 Hexachlorobutane	500U	118-74-1 Hexachlorobenzene - U J	500U
50-95-3 Nitrobenzene	2000U	87-86-5 Pentachlorophenol	500U
50-91-1 Isophorone	800U	85-01-8 Phenanthrene	300U
88-75-5 2-Nitrophenol	500U	120-12-7 Anthracene - U J	400U
105-67-9 2,4-Dimethylphenol	300U	84-74-2 Di-n-Butylphthalate - J S	700JB
85-0 Benzoic Acid U J	1000U	206-44-0 Fluoranthene	600U
1-91-1 bis (2-Chloroethoxy) Methane	900U	92-87-5 Benzidine - R	2000U
120-83-2 2,4-Dichlorophenol	300U	129-00-0 Pyrene	2000U
0-82-1 1,2,4-Trichlorobenzene	600U	85-68-7 Butylbenzylphthalate	800U
-20-3 Naphthalene	600U	91-94-1 3,3'-Dichlorobenzidine - R	1000U
106-47-8 4-Chloraniline U J	3000U	56-55-3 Benzo(a)anthracene	400U
77-68-3 Hexachlorobutadiene	800U	117-81-7 bis (2-Ethylhexyl) Phthalate J	1100B
50-7 4-Chloro-3-Methylphenol	500U	218-01-9 Chrysene	400U
91-57-6 2-Methylnaphthalene	300U	117-84-0 Di-n-octyl Phthalate	400U
77-47-4 Hexachlorocyclopentadiene	2000U	205-99-2 Benzo(b)Fluoranthene	1000U
50-2 2,4,6-Trichlorophenol	300U	207-08-9 Benzo(k)Fluoranthene U J	500U
50-4 2,4,5-Trichlorophenol	300U	50-32-8 Benzo(a)Pyrene	400U
91-58-7 2-Chloronaphthalene	500U	193-39-5 Indeno(1,2,3-cd)Pyrene	400U
50-4 2-Nitroaniline - R	2000U	53-70-3 Dibenzo(a,h)Anthracene	400U
11-11-3 Dimethyl Phthalate	500U	191-24-2 Benzo(g,h,i)Perylene	500U
208-96-8 Acenaphthylene - U J	300U		
90-09-2 3-Nitroaniline - U J	2000U	(1)- Cannot be separated from diphenylamine	

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Sample Number

H1843

28
10/27/85

Organics Analysis Data Sheet
(Page 3)

ANALYSIS NOT REQUIRED

Pesticide/PCBs

Concentration: Low Medium (Circle One)

Date Extracted/Prepared: _____

Date Analyzed: _____

Conc/Dil Factor: _____

CAS Number	ug/l or ug/Kg (Circle One)
519-84-6	Alpha-BHC
519-85-7	Beta-BHC
519-86-8	Delta-BHC
58-89-9	Gamma-BHC (Lindane)
76-44-8	Heptachlor
309-00-2	Aldrin
1024-57-3	Heptachlor Epoxide
959-98-8	Endosulfan I
60-57-1	Dieldrin
72-55-9	4, 4'-DDE
72-20-8	Endrin
33213-65-9	Endosulfan II
72-54-8	4, 4'-DDD
7421-93-4	Endrin Aldehyde
1031-07-8	Endosulfan Sulfate
50-29-3	4, 4'-DDT
72-43-5	Methoxychlor
53494-70-5	Endrin Ketone
57-74-9	Chlordane
8001-35-2	Tetraphene
12674-11-2	Aroclor-1016
11104-28-2	Aroclor-1221
11141-16-5	Aroclor-1232
53469-21-9	Aroclor-1242
12672-29-6	Aroclor-1248
11097-69-1	Aroclor-1254
11096-82-5	Aroclor-1260

V_i = Volume of extract injected (ul)

V_s = Volume of water extracted (ml)

W_s = Weight of sample extracted (g)

V_t = Volume of total extract (ul)

V_s _____ or W_s _____ V_t _____ V_i _____

Environmental Protection Agency, CLP Sample Management Office,
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Sample Number
H1843

Organics Analysis Data Sheet
(Page 4)

Tentatively Identified Compounds

No volatile compounds found.

GAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or ug/kg)
1. 96-48-0	2(3H)-Furanone, dihydro-	BNA	599	100J
2. 628-68-2	Ethanol, 2,2'-oxybis-, diacetate	BNA	704	1000J
3. --	Unknown alcohol	BNA	723	90J
4. --	Unknown	BNA	729	300J
5. --	Unknown	BNA	1010	200J
6. --	Unknown	BNA	1093	700J
7. --	Unknown	BNA	1287	1000J
8. --	Unknown fatty acid	BNA	1471	100J
9. 10544-50-0	Sulfur, mol. (S8)	BNA	1519	100000J
10. --	Unknown amide	BNA	1712	700J
11. --	Unknown	BNA	1773	200J
12. --	Unknown	BNA	1860	300J
13. --	Unknown saturated aliphatic hydrocarbon	BNA	1871	200J
14. --	Unknown saturated aliphatic hydrocarbon	BNA	1976	600J
15. --	Unknown	BNA	2062	500J
16. --	Unknown saturated aliphatic hydrocarbon	BNA	2116	300J
17. --	Unknown	BNA	2127	200J
18. 57-88-5	Cholesterol	BNA	2205	1000J
19. 83-47-6	Stigmast-5-en-3-ol, (3, beta., 24S)-	BNA	2427	500J
20.				
21.				
22.				
23.				
24.				
25.				
26.				
27.				
28.				
29.				
30.				

26
ppm

Organics Analysis Data Sheet
(Page 1)

Laboratory Name: CE-EMSI

Lab Sample ID No: 043085V13 (VOA) 052285C12 (EPA)

Sample Matrix: Soil

Release Authorized By: Richard J Coutas

Case No: 4242

QC Report No: 4242

Contract No: 68-01-6783

Date Sample Received: 24-APR-85

Volatile Compounds

Concentration: Low
Date Extracted/Prepared: NA
Date Analyzed: 30-APR-85
Conc/Dil Factor: NA pH 8.3
Percent Moisture: 14.53
Percent Moisture (Decanted): NA

IS Number	ug/l or ug/kg (Circle One)	CAS Number	ug/l or ug/kg (Circle One)
-87-3	Chloromethane -	30U	79-34-5 1,1,2,2-Tetrachloroethane <u>J</u>
74-83-9	Bromomethane	2U	78-87-5 1,2-Dichloropropane <u>J</u>
75-01-4	Vinyl Chloride <u>J</u>	30U	10061-02-6 Trans-1,3-Dichloropropene
-00-3	Chloroethane	2U	79-01-6 Trichloroethene
-09-2	Methylene Chloride <u>J</u>	15B	124-48-1 Dibromochloromethane
67-64-1	Acetone <u>J</u>	3B	79-00-5 1,1,2-Trichloroethane
-74-15-0	Carbon Disulfide	5U	71-43-2 Benzene
-35-4	1,1-Dichloroethene	3U	10061-01-5 cis-1,3-Dichloropropene
75-34-3	1,1-Dichloroethane	2U	110-75-8 2-Chloroethylvinylether <u>R</u>
156-60-5	Trans-1,2-Dichloroethene	2U	75-25-2 Bromoform
-66-3	Chloroform	1U	591-78-6 2-Hexanone <u>J</u>
207-06-2	1,2-Dichloroethane	1U	108-10-1 4-Methyl-2-Pentanone
78-93-3	2-Butanone <u>J</u>	1U	127-18-4 Tetrachloroethene
-55-6	1,1,1-Trichloroethane <u>J</u>	3U	108-88-3 Toluene - <u>J</u>
-23-5	Carbon Tetrachloride	3U	108-90-7 Chlorobenzene
108-05-4	Vinyl Acetate <u>R</u>	2U	100-41-4 Ethylbenzene
75-27-4	Bromodichloromethane	1U	100-42-5 Styrene
			Total Xylenes

Data Reporting Qualifiers

For reporting results to EPA, the following results qualifiers are used.

G If the result is a value greater than or equal to the detection limit, report the value

G This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides >10 ng/uL in the final extract should be confirmed by GC/MS.

B This flag is used when the analyte is found in the blank as well as a sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action.

S Indicates compound was spiked into the sample for the matrix recovery study. The amounts that were spiked into and recovered from the sample are reported on Form III.

3 187

4/84

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Sample Number
H1844

72
shelby

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Organics Analysis Data Sheet
(Page 2)

Semi-volatile Compounds

Concentration: Low
Date Extracted/Prepared: 30-APR-85
Date Analyzed: 22-MAY-85 ✓
Conc/Dil Factor: 1.0

CAS Number	ug/l or $\mu\text{g}/\text{kg}$ (Circle One)	CAS Number	ug/l or $\mu\text{g}/\text{kg}$ (Circle One)		
62-75-9	N-Nitrosodimethylamine	500U	83-32-9	Acenaphthene	300U
10-95-2	Phenol J8	200U	51-28-5	2,4-Dinitrophenol ~ U J	2000U
12-53-3	Aniline U J	300U	100-02-7	4-Nitrophenol ~ U J	2000U
111-44-4	Bis(2-Chloroethyl)Ether U J	300U	132-64-9	Dibenzofuran	200U
55-57-8	2-Chlorophenol	200U	121-14-2	2,4-Dinitrotoluene	600U
41-73-1	1,3-Dichlorobenzene	200U	606-20-2	2,6-Dinitrotoluene	500U
106-46-7	1,4-Dichlorobenzene	200U	84-66-2	Diethylphthalate	300U
100-51-6	Benzyl Alcohol	500U	7005-72-3	4-Chlorophenyl-phenylether	1000U
5-50-1	1,2-Dichlorobenzene	200U	86-73-7	Fluorene	200U
55-48-7	2-Methylphenol	200U	100-01-6	4-Nitroaniline - U J	600U
39538-32-9	Bis (2-Chloroisopropyl) Ether	200U	534-52-1	4,6-Dinitro-2-Methylphenol	1000U
56-44-5	4-Methylphenol	800U	86-30-6	N-Nitrosodiphenylamine (1)	500U
21-64-7	N-Nitroso-Di-n-Propylamine	600U	101-55-3	4-Bromophenylphenylether - U J	400U
67-72-1	Benzachloroethene	200U	118-74-1	Benzachlorobenzene - U J	200U
99-95-3	Nitrobenzene	600U	87-86-5	Pentachlorophenol	200U
8-59-1	Lophorone	300U	85-01-8	Phenanthrene	100U
55-75-5	2-Nitrophenol	200U	120-12-7	Anthracene - U J	200U
105-67-9	2,4-Dinethylphenol	100U	84-74-2	Di-n-Butylphthalate - J8	540UB
5-85-0	Benzoic Acid U J	500U	206-44-0	Fluoranthene	300U
111-91-1	Bis (2-Chloroethoxy) Methane	400U	92-87-5	Benzidine - R	800U
120-83-2	2,4-Dichlorophenol	100U	129-00-0	Pyrene	600U
20-82-1	1,2,4-Trichlorobenzene	300U	85-68-7	Butylbenzylphthalate	300U
1-20-3	Naphthalene	300U	91-94-1	3,3'-Dichlorobenzidine - R	500U
106-47-8	4-Chloroaniline U J	1000U	56-55-3	Benzo(a)anthracene	200U
97-68-3	Benzachlorobutadiene	300U	117-81-7	Bis (2-Ethylhexyl) Phthalate - J	290B
9-50-7	4-Chloro-3-Methylphenol	200U	218-01-9	Chrysene	200U
51-57-6	2-Methylnaphthalene	100U	117-84-0	Di-n-octyl Phthalate	200U
77-47-4	Benzachlorocyclopentadiene	800U	205-99-2	Benzo(b)Fluoranthene	600U
8-06-2	2,4,6-Trichlorophenol	100U	207-08-9	Benzo(k)Fluoranthene U J	200U
5-95-4	2,4,5-Trichlorophenol	100U	50-32-8	Benzo(a)Pyrene	200U
91-58-7	2-Chloronaphthalene	200U	193-39-5	Indeno(1,2,3-cd)Pyrene	200U
78-74-4	2-Nitroaniline - R	1000U	53-70-3	Dibenzo(a,h)Anthracene	200U
31-11-3	Dimethyl Phthalate	200U	191-24-2	Benzo(g,h,i)Perylene	200U
208-96-8	Acenaphthylene - U J	100U			
99-09-2	3-Nitroaniline - U J	1000U	(1)- Cannot be separated from diphenylamine		

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Sample Number
H1844

12/13
12/15r

Organics Analysis Data Sheet
(Page 3)

ANALYSIS NOT REQUIRED

Pesticide/PCBs

Concentration: Low Medium (Circle One)

Date Extracted/Prepared: _____

Date Analyzed: _____

Conc/Dil Factor: _____

CAS Number		ug/l or ug/Kg (Circle One)
319-84-6	Alpha-BHC	
319-85-7	Beta-BHC	
319-86-8	Delta-BHC	
58-89-9	Gamma-BHC (Lindane)	
76-44-8	Heptachlor	
309-00-2	Aldrin	
1024-57-3	Heptachlor Epoxide	
959-98-8	Endosulfan I	
60-57-1	Dieldrin	
72-55-9	4, 4'-DDE	
72-20-8	Endrin	
33213-65-9	Endosulfan II	
72-54-8	4, 4'-DDD	
7421-93-4	Endrin Aldehyde	
1031-07-8	Endosulfan Sulfate	
50-29-3	4, 4'-DDT	
72-43-5	Methoxychlor	
53494-70-5	Endrin Ketone	
57-74-9	Chlordane	
8001-35-2	Toxaphene	
12674-11-2	Aroclor-1016	
11104-28-2	Aroclor-1221	
11141-16-5	Aroclor-1232	
53469-21-9	Aroclor-1242	
12672-29-6	Aroclor-1248	
11097-69-1	Aroclor-1254	
11096-82-5	Aroclor-1260	

V_i = Volume of extract injected (ul)

V_s = Volume of water extracted (ml)

W_s = Weight of sample extracted (g)

V_t = Volume of total extract (ul)

V_s _____ or W_s _____ V_t _____ V_i _____

Sample Number

R1844

12
4/78

Organics Analysis Data Sheet
(Page 4)

Tentatively Identified Compounds

No volatile compounds found.

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or mg/kg)
1. 96-48-0	2(3H)-Furanone, dihydro-	BNA	586	200JB
2. --	Unknown	BNA	589	200J
3. 628-68-2	Ethanol, 2,2'-oxobis-, diacetate	BNA	701	1000JB
4. 104-76-7	1-Hexanol, 2-ethyl-	BNA	719	200J
5. 57-10-3	Palmitic acid	BNA	1477	200J
6. --	Unknown amide	BNA	1719	700J
7. 103-23-1	Hexanedioic acid, bis(2-ethylhexyl)ester	BNA	1732	300JB
8. --	Unknown	BNA	1780	200J
9. --	Unknown saturated aliphatic hydrocarbon	BNA	1877	400J
10. --	Unknown saturated aliphatic hydrocarbon	BNA	1982	1000J
11. --	Unknown	BNA	2083	300J
12. --	Unknown saturated aliphatic hydrocarbon	BNA	2122	900J
13. --	Unknown	BNA	2132	400J
14. --				
15. --				
16. --				
17. --				
18. --				
19. --				
20. --				
21. --				
22. --				
23. --				
24. --				
25. --				
26. --				
27. --				
28. --				
29. --				
30. --				

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Sample Number
RI 845

Re
Dolan

Organics Analysis Data Sheet
(Page 1)

Laboratory Name: CE-DMSI
Lab Sample ID No: 05018SV09 (VOA) 05228SC15 (EPA)
Sample Matrix: Soil
Data Release Authorized By: Richard Lourdes

Case No: 4242
QC Report No: 4242
Contract No: 68-01-6783
Date Sample Received: 24-APR-85

Volatile Compounds

Concentration: Low
Date Extracted/Prepared: NA
Date Analyzed: 01-MAY-85
Conc/Dil Factor: NA pH 7.7
Percent Moisture: 28.32
Percent Moisture (Decanted): NA

S Number	ug/l or ug/Kg (Circle one)	CAS Number	ug/l or ug/Kg (Circle one)
1-87-3	Chloromethane	300	79-34-5 1,1,2,2-Tetrachloroethane W J
1-83-9	Bromomethane	30	78-87-5 1,2-Dichloropropane W J
75-01-4	Vinyl Chloride W J	300	10061-02-6 Trans-1,3-Dichloropropene
73-00-3	Chloroethane	30	79-01-6 Trichloroethene
5-09-2	Methylene Chloride J	40B	124-48-1 Dibromochloromethane
67-64-1	Acetone J	33B	79-00-5 1,1,2-Trichloroethane
75-15-0	Carbon Disulfide	60	71-43-2 Benzene
5-35-4	1,1-Dichloroethene	40	10061-01-5 cis-1,3-Dichloropropene
5-34-3	1,1-Dichloroethane	30	110-75-8 2-Chloroethylvinylether R
156-60-5	Trans-1,2-Dichloroethene	30	75-25-2 Bromoform
7-66-3	Chloroform	100	591-78-6 2-Hexanone W J
37-06-2	1,2-Dichloroethane	100	108-10-1 4-Methyl-2-Pentanone
78-93-3	2-Butanone T	5B	127-18-4 Tetrachloroethene
71-55-6	1,1,1-Trichloroethane W J	40	108-88-3 Toluene J B
5-23-5	Carbon Tetrachloride	40	108-90-7 Chlorobenzene
108-05-4	Vinyl Acetate R	30	100-41-4 Ethylbenzene
75-27-4	Bromodichloromethane	100	100-42-5 Styrene
			Total Xylenes

Data Reporting Qualifiers

For reporting results to EPA, the following results qualifiers are used.

V Value: If the result is a value greater than or equal to the detection limit, report the value

C This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides ≥ 10 ng/ul in the final extract should be confirmed by GC/MS.

U Compound was analyzed for but not detected. The number is the minimum attainable detection limit for the sample.

B This flag is used when the analyte is found in the blank as well as a sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action.

J Indicates an estimated value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicates the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero (e.g. 10J).

S Indicates compound was spiked into the sample for the matrix recovery study. The amounts that were spiked into and recovered from the sample are reported on Form III.

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Sample Number
H1845

2/1/85

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Organics Analysis Data Sheet
(Page 2)

Semivolatile Compounds

Concentration: Low
Date Extracted/Prepared: 30-APR-85
Date Analyzed: 22-MAY-85
Conc/Dil Factor: 2.0

CAS Number	ug/l or $\mu\text{g}/\text{kg}$ (Circle One)	CAS Number	ug/l or $\mu\text{g}/\text{kg}$ (Circle One)		
62-75-9	N-Nitrosodimethylamine	1000U	83-32-9	Acenaphthene	500U
108-95-2	Phenol β	300U	51-28-5	2,4-Dinitrophenol - α , β	4000U
111-53-3	Aniline α , β	600U	100-02-7	4-Nitrophenol - α , β	4000U
111-44-4	bis(2-Chloroethyl)Ether α , β	600U	132-64-9	Dibenzo furan	400U
95-57-8	2-Chlorophenol	300U	121-14-2	2,4-Dinitrotoluene	1000U
11-73-1	1,3-Dichlorobenzene	300U	606-20-2	2,6-Dinitrotoluene	1000U
126-46-7	1,4-Dichlorobenzene	300U	84-66-2	Diethylphthalate	500U
100-51-6	Benzyl Alcohol	1000U	7005-72-3	4-Chlorophenyl-phenylether	200U
11-50-1	1,2-Dichlorobenzene	300U	86-73-7	Fluorene	300U
11-48-7	2-Methylphenol	300U	100-01-6	4-Nitroaniline - α , β	1000U
39638-32-9	bis (2-Chloroisopropyl) Ether	300U	534-52-1	4,6-Dinitro-2-Methylphenol	3000U
106-44-5	4-Methylphenol	2000U	86-30-6	N-Nitrosodiphenylamine (1)	1000U
11-64-7	N-Nitroso-Di-n-Propylamine	1000U	101-55-3	4-Bromophenylphenylether - α , β	700U
67-72-1	Hexachloroethane	400U	118-74-1	Hexachlorobenzene - α , β	400U
98-95-3	Nicrobenzene	1000U	87-86-5	Pentachlorophenol	400U
3-59-1	Iso phorone	600U	85-01-8	Phenanthrene	200U
1-375-5	2-Nitrophenol	400U	120-12-7	Anthracene - α , β	300U
105-67-9	2,4-Dimethylphenol	200U	84-74-2	Di-n-Butylphthalate - β	660U B
11-85-0	Benzoic Acid α , β	1000U	206-44-0	Fluoranthene	500U
11-91-1	bis (2-Chloroethoxy) Methane	700U	92-87-5	Benzidine - R	2000U
120-83-2	2,4-Dichlorophenol	200U	129-00-0	Pyrene	1000U
120-82-1	1,2,4-Trichlorobenzene	500U	85-66-7	Butylbenzylphthalate	600U
1-20-3	Naphthalene	500U	91-94-1	3,3'-Dichlorobenzidine - R	1000U
106-47-8	4-Chloroaniline α , β	3000U	56-55-3	Benzo(a)anthracene	300U
87-68-3	Hexachlorobutadiene	600U	117-61-7	bis (2-Ethylhexyl) Phthalate β	840B
11-50-7	4-Chloro-3-Methylphenol	400U	218-01-9	Chrysene	300U
1-157-6	2-Methylnaphthalene	200U	117-84-0	Di-n-octyl Phthalate	300U
77-47-4	Hexachlorocyclopentadiene	2000U	205-99-2	Benzo(b)Fluoranthene	1000U
78-06-2	2,4,6-Trichlorophenol	200U	207-08-9	Benzo(k)Fluoranthene β	400U
5-95-4	2,4,5-Trichlorophenol	200U	50-32-8	Benzo(a)Pyrene	300U
91-58-7	2-Chloronaphthalene	400U	193-39-5	Indeno(1,2,3-cd)Pyrene	300U
88-74-4	2-Nitroaniline - R	2000U	53-70-3	Dibenzo(a,h)Anthracene	300U
31-11-3	Dimethyl Phthalate	400U	191-24-2	Benzo(g,h,i)Perylene	400U
708-96-8	Acenaphthylene - α , β	200U			
99-09-2	3-Nitroaniline - α , β	2000U			
			(1) Cannot be separated from diphenylamine		

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Sample Number
H1845

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Organics Analysis Data Sheet
(Page 3)

ANALYSIS NOT REQUIRED

Pesticide/PCBs

Concentration: Low Medium (Circle One)

Date Extracted/Prepared: _____

Date Analyzed: _____

Conc/Dil Factor: _____

CAS Number		ug/l or ug/Kg (Circle One)
319-84-6	Alpha-BHC	
319-85-7	Beta-BHC	
319-86-8	Delta-BHC	
58-89-9	Gamma-BHC (Lindane)	
76-44-8	Heptachlor	
309-00-2	Aldrin	
1024-57-3	Heptachlor Epoxide	
1959-98-8	Endosulfan I	
60-57-1	Dieldrin	
72-55-9	4, 4'-DDE	
72-20-8	Endrin	
33213-65-9	Endosulfan II	
72-54-8	4, 4'-DDD	
7421-93-4	Endrin Aldehyde	
1031-07-8	Endosulfan Sulfate	
50-29-3	4, 4'-DDT	
72-43-5	Methoxychlor	
53494-70-5	Endrin Ketone	
57-74-9	Chlordane	
8001-35-2	Toxaphene	
12674-11-2	Aroclor-1016	
11104-28-2	Aroclor-1221	
11141-16-5	Aroclor-1232	
53459-21-9	Aroclor-1242	
12672-29-6	Aroclor-1248	
11097-69-1	Aroclor-1254	
11096-82-5	Aroclor-1260	

V_i = Volume of extract injected (uL)

V_s = Volume of water extracted (mL)

W_s = Weight of sample extracted (g)

V_t = Volume of total extract (uL)

V_s _____ or W_s _____ V_t _____ V_i _____

Sample Number
H1845

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4/19/86

Organics Analysis Data Sheet
(Page 4)

Tentatively Identified Compounds

No volatile compounds found.

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or kg/kg)
1. 96-48-0	2(3H)-Furanone, dihydro-	BNA	581	200JB
2. 628-68-2	Ethanol, 2,2'-oxybis-, diacetate	BNA	694	2000JB
3. 104-76-7	1-Hexanol, 2-ethyl-	BNA	714	300J
4. --	Unknown	BNA	720	500J
5. --	Unknown	BNA	1009	700J
6. --	Unknown	BNA	1092	2000J
7. --	Unknown	BNA	1286	1000J
8. 17233-71-5	Hexathiepane	BNA	1352	500J
9. 10544-50-0	Sulfur, mol. (S8)	BNA	1512	100000J
10. --	Unknown amide	BNA	1712	1000J
11. 103-23-1	Hexanedioic acid, bis(2-ethylhexyl)ester	BNA	1725	300JB
12. --	Unknown saturated aliphatic hydrocarbon	BNA	1975	800J
13. 57-88-5	Cholesterol	BNA	2205	1000J
14. --	Unknown sterol	BNA	2426	700J
15. --				
16. --				
17. --				
18. --				
19. --				
20. --				
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27. --				
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Sample Number
H1846

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MS-16
being filmed

Organics Analysis Data Sheet
(Page 1)

Laboratory Name: CE-EMSI

Sample ID No: 043085V15 (VDA) 052385C10 (ENA)

Sample Matrix: Soil

Data Release Authorized By:

Richard Klontz

Case No: 4242

QC Report No: 4242

Contract No: 68-01-6783

Date Sample Received: 24-APR-85

Volatile Compounds

Concentration: Low
Date Extracted/Prepared: NA
Date Analyzed: 30-APR-85
Conc/Dil Factor: NA pH 7.95
Percent Moisture: 6.78
Percent Moisture (Decanted): NA

I S number	ug/l or ug/kg (Circle one)	CAS Number	ug/l or ug/kg (Circle one)
-87-3	Chloromethane	300	79-34-5 1,1,2,2-Tetrachloroethane WJ
-83-9	Bromomethane	20	78-87-5 1,2-Dichloropropane WJ
75-01-4	Vinyl Chloride WJ	300	10061-02-6 Trans-1,3-Dichloropropene
75-00-3	Chloroethane	20	79-01-6 Trichloroethane
-09-2	Methylene Chloride J	150	124-48-1 Dibromochloromethane
67-64-1	Acetone J	100	79-00-5 1,1,2-Trichloroethane
75-15-0	Carbon Disulfide	40	71-43-2 Benzene
-35-4	1,1-Dichloroethene	30	10061-01-5 cis-1,3-Dichloropropene
-34-3	1,1-Dichloroethane	20	110-75-8 2-Chloroethylvinylether R
156-60-5	Trans-1,2-Dichloroethene	20	75-25-2 Bromoform
-66-3	Chloroform	100	591-78-6 2-Hexanone WJ
17-06-2	1,2-Dichloroethane	100	106-10-1 4-Methyl-2-Pentanone
78-93-3	2-Butanone R	100	127-18-4 Tetrachloroethene
71-55-6	1,1,1-Trichloroethane WJ	30	106-88-3 Toluene
-23-5	Carbon Tetrachloride	30	108-90-7 Chlorobenzene
106-05-4	Vinyl Acetate R	20	100-41-4 Ethylbenzene
75-27-4	Bromodichloromethane	100	100-42-5 Styrene
			Total Xylenes

Data Reporting Qualifiers

For reporting results to EPA, the following results qualifiers are used.

V Value If the result is a value greater than or equal to the detection limit, report the value

U Compound was analyzed for but not detected.
The number is the minimum attainable detection limit for the sample.

J Indicates an estimated value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicates the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero (e.g. 10).

C This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides >10 ng/uL in the final extract should be confirmed by GC/MS.

B This flag is used when the analyte is found in the blank as well as a sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action.

S Indicates compound was spiked into the sample for the matrix recovery study. The amounts that were spiked into and recovered from the sample are reported on Form III.

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Sample Number
EN846

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4/16/85

Organics Analysis Data Sheet
(Page 2)

Semivolatile Compounds

Concentration: Low
Date Extracted/Prepared: 30-APR-85
Date Analyzed: 23-MAY-85
Conc/Dil Factor: 10.0

CAS Number	ug/l or ug/kg (Circle one)	CAS Number	ug/l or ug/kg (Circle one)		
62-75-9	N-Nitrosodimethylamine	5000	83-32-9	Acenaphthene	3000
108-95-2	Phenol Jβ	2000	51-28-5	2,4-Dinitrophenol - u J	20000
115-3-3	Aniline u J	3000	100-02-7	4-Nitrophenol - u J	20000
111-44-4	bis(2-Chloroethyl)Ether u J	3000	132-64-9	Dibenzofuran	2000
95-57-8	2-Chlorophenol	2000	121-14-2	2,4-Dinitrotoluene	6000
111-73-1	1,3-Dichlorobenzene	2000	606-20-2	2,6-Dinitrotoluene	5000
106-46-7	1,4-Dichlorobenzene	2000	84-66-2	Diethylphthalate	3000
100-51-6	Benzyl Alcohol	5000	7005-72-3	4-Chlorophenyl-phenylether	1000
111-50-1	1,2-Dichlorobenzene	2000	86-73-7	Fluorene	2000
111-48-7	2-Methylphenol	2000	100-01-6	4-Nitroaniline - u J	6000
39638-32-9	bis (2-Chloroisopropyl) Ether	2000	534-52-1	4,6-Dinitro-2-Methylphenol	10000
106-44-5	4-Methylphenol	8000	86-30-6	N-Nitrosodiphenylamine (1)	5000
111-64-7	N-Nitroso-Di-n-propylamine	6000	101-55-3	4-Bromophenylphenylether - u J	4000
111-72-1	Hexachloroethane	2000	118-74-1	Hexachlorobenzene - u J	2000
98-95-3	Nitrobenzene	6000	87-86-5	Pentachlorophenol	2000
111-59-1	Isophorone	3000	85-01-8	Phenanthrene	1000
111-75-5	2-Nitrophenol	2000	120-12-7	Anthracene - u J	2000
105-67-9	2,4-Dimethylphenol	1000	84-74-2	Di-n-butylphthalate Jβ	10000
111-85-0	Benzoic Acid u J	5000	206-44-0	Fluoranthene	3000
111-91-1	bis (2-Chloroethyl) Methane	4000	92-87-5	Benzidine - R	8000
120-83-2	2,4-Dichlorophenol	1000	129-00-0	Pyrene	6000
120-82-1	1,2,4-Trichlorobenzene	3000	85-68-7	Ethylbenzylphthalate	3000
111-20-3	Naphthalene	3000	91-94-1	3,3'-Dichlorobenzidine - R	5000
111-06-47-8	4-Chloroaniline - u J	10000	56-55-3	Benzo(a)anthracene	2000
87-68-3	Hexachlorobutadiene	3000	117-81-7	bis (2-Ethylhexyl) Phthalate R	2000
98-50-7	4-Chloro-3-Methylphenol	2000	218-01-9	Chrysene	2000
111-57-6	2-Methylnaphthalene	1000	117-84-0	Di-n-octyl Phthalate	2000
77-47-4	Hexachlorocyclopentadiene	8000	205-99-2	Benzo(b)Fluoranthene	6000
98-06-2	2,4,6-Trichlorophenol	1000	207-08-9	Benzo(k)Fluoranthene u J	2000
5-95-4	2,4,5-Trichlorophenol	1000	50-32-8	Benzo(a)Pyrene	2000
91-58-7	2-Chloronaphthalene	2000	193-39-5	Indeno(1,2,3-cd)Pyrene	2000
88-74-4	2-Nitroaniline - R	10000	53-70-3	Dibenzo(a,h)Anthracene	2000
31-11-3	Dimethyl Phthalate	2000	191-24-2	Benzo(g,h,i)Perylene	2000
111-08-8	Acenaphthylene - u J	1000			
99-09-2	3-Nitroaniline - u J	10000	(1)- Cannot be separated from diphenylamine		

Sample Number

H1846

26
ref/hs

Organics Analysis Data Sheet
(Page 3)

ANALYSIS NOT REQUIRED

Pesticide/PCBs

Concentration: Low Medium (Circle One)

Date Extracted/Prepared: _____

Date Analyzed: _____

Conc/Dil Factor: _____

CAS Number		ug/l or ug/Kg (Circle One)
319-84-6	Alpha-BHC	
319-85-7	Beta-BHC	
319-86-8	Delta-BHC	
58-89-9	Gamma-BHC (Lindane)	
76-44-8	Heptachlor	
309-00-2	Aldrin	
1024-57-3	Heptachlor Epoxide	
959-96-8	Endosulfan I	
60-57-1	Dieldrin	
72-55-9	4, 4'-DDE	
72-20-8	Endrin	
33213-65-9	Endosulfan II	
72-54-8	4, 4'-DDD	
7421-93-4	Endrin Aldehyde	
1031-07-8	Endosulfan Sulfate	
50-29-3	4, 4'-DDT	
72-43-5	Methoxychlor	
53494-70-5	Endrin Ketone	
57-74-9	Chlordane	
8001-35-2	Toxaphene	
12674-11-2	Aroclor-1016	
11104-28-2	Aroclor-1221	
11141-16-5	Aroclor-1232	
53469-21-9	Aroclor-1242	
12672-29-6	Aroclor-1248	
11097-69-1	Aroclor-1254	
11096-82-5	Aroclor-1260	

V_i = Volume of extract injected (ul)

V_s = Volume of water extracted (ml)

W_s = Weight of sample extracted (g)

V_t = Volume of total extract (ul)

V_s _____ or W_s _____ V_t _____ V_i _____

3 266

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Sample Number:

H1846

78
relabel

Organics Analysis Data Sheet
(Page 4)

Tentatively Identified Compounds

No volatile compounds found.

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or ug/kg)
1. 96-48-0	2(3H)-Furanone, dihydro-	BNA	578	700JB
2. 542-10-9	1,1-Ethanediol, diacetate	BNA	688	1000J
3. --	Unknown	BNA	1443	2000J
4. --	Unknown	BNA	1691	3000J
5. --	Unknown	BNA	2144	3000J
6.				
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10 Box 816. Alexandria, Virginia 22313 703/557-2490

Sample Number
H1847

Organics Analysis Data Sheet
(Page 1)

Laboratory Name: CE-PMI

Sample ID No: 043065V14 (VDA) 052285CL3 (ENA)

Sample Matrix: Soil

Data Release Authorized By:

Richard J Coontess

Case No: 4242

QC Report No: 4242

Contract No: 68-01-6783

Date Sample Received: 24-APR-85

Volatile Compounds

Concentration: Low
Date Extracted/Prepared: NA
Date Analyzed: 30-APR-85
Conc/Dil Factor: NA pH 8.3
Percent Moisture: 4.15
Percent Moisture (Decanted): NA

I S Number		ug/l or ug/Kg (Circle one)	CAS Number		ug/l or ug/Kg (Circle one)
-87-3	Chloromethane	2U	79-34-5	1,1,2,2-Tetrachloroethane UJ	X
-83-9	Bromomethane	2U	78-87-5	1,2-Dichloropropane UJ	X
75-01-4	Vinyl Chloride	2U	10061-02-6	Trans-1,3-Dichloropropene	X
75-00-3	Chloroethane	2U	79-01-6	Trichloroethene	X
-09-2	Methylene Chloride J	10B	124-48-1	Dibromochloromethane	X
67-64-1	Acetone J	3B	79-00-5	1,1,2-Trichloroethane	X
75-15-0	Carbon Disulfide	4U	71-43-2	Benzene	X
-35-4	1,1-Dichloroethene	3U	10061-01-5	cis-1,3-Dichloropropene	X
-75-34-3	1,1-Dichloroethane	2U	110-75-8	2-Chloroethylvinylether R	X
156-60-5	Trans-1,2-Dichloroethene	2U	75-25-2	Bromoform	X
-66-3	Chloroform	1U	591-78-6	2-Hexanone UJ	X
77-06-2	1,2-Dichloroethane	1U	108-10-1	4-Methyl-2-Pentanone	X
78-93-3	2-Butanone J	1U	127-18-4	Tetrachloroethene	X
71-55-6	1,1,1-Trichloroethane UJ	3U	108-88-3	Toluene U-B	7B
-23-5	Carbon Tetrachloride	3U	108-90-7	Chlorobenzene	X
106-05-4	Vinyl Acetate R	2U	100-41-4	Ethylbenzene	X
75-27-4	Bromodichloromethane	1U	100-42-5	Styrene	X
				Total Xylenes	X

Data Reporting Qualifiers

For reporting results to EPA, the following results qualifiers are used.

V Value If the result is a value greater than or equal to the detection limit, report the value

U Compound was analyzed for but not detected. The number is the minimum attainable detection limit for the sample.

J Indicates an estimated value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicates the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero (e.g. 10J).

C This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides >10 ng/uL in the final extract should be confirmed by GC/MS.

B This flag is used when the analyte is found in the blank as well as a sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action.

S Indicates compound was spiked into the sample for the matrix recovery study. The amounts that were spiked into and recovered from the sample are reported on Form III.

6/9/85

Organics Analysis Data Sheet
(Page 2)

Semivolatile Compounds

Concentration: Low
Date Extracted/Prepared: 30-APR-85
Date Analyzed: 22-MAY-85
Conc/Dil Factor: 1.0

CAS Number	ug/l or ug/kg (Circle one)	CAS Number	ug/l or ug/kg (Circle one)
2-75-9 N-Nitrosodimethylamine	500U	83-32-9 Acenaphthene	300U
108-95-2 Phenol - J S	200U	51-28-5 2,4-Dinitrophenol - U J	2000U
2-53-3 Aniline - U J	300U	100-02-7 4-Nitrophenol - U J	2000U
11-44-4 bis(2-Chloroethyl)Ether U J	300U	132-64-9 Dibenzofuran	200U
95-17-8 2-Chlorophenol	200U	121-14-2 2,4-Dinitrotoluene	700U
54-17-1 1,3-Dichlorobenzene	200U	606-20-2 2,6-Dinitrotoluene	500U
66-46-7 1,4-Dichlorobenzene	200U	84-66-2 Diethylphthalate	300U
100-51-6 Benzyl Alcohol	500U	7005-72-3 4-Chlorophenyl-phenylether	100U
95-50-1 1,2-Dichlorobenzene	200U	86-73-7 Fluorene	200U
54-48-7 2-Methylphenol	200U	100-01-6 4-Nitroaniline - U J	700U
9638-32-9 bis (2-Chloroisopropyl) Ether	200U	534-52-1 4,6-Dinitro-2-Methylphenol	1000U
106-44-5 4-Methylphenol	800U	N-Nitroso-diphenylamine (1)	500U
21-64-7 N-Nitroso-Di- α -Propylamine	700U	4-Bromophenylphenylether - U J	400U
7-72-1 Hexachloroethane	200U	118-74-1 Hexachlorobenzene - U J	200U
98-95-3 Nitrobenzene	700U	87-86-5 Pentachlorophenol	200U
78-59-1 Isophorone	300U	85-01-8 Phenanthrene	100U
108-75-5 2-Nitrophenol	200U	120-12-7 Anthracene - U J	200U
120-67-9 2,4-Dimethylphenol	100U	84-74-2 Di-n-Butylphthalate - J S	450JB
65-85-0 Benzoic Acid U J	500U	206-44-0 Fluoranthene	300U
11-91-1 bis (2-Chloroethoxy) Methane	400U	92-87-5 Benzidine - R	300U
20-83-2 2,4-Dichlorophenol	100U	129-00-0 Pyrene	700U
120-82-1 1,2,4-Trichlorobenzene	300U	85-68-7 Butylbenzylphthalate	300U
71-20-3 Naphthalene	300U	91-94-1 3,3'-Dichlorobenzidine - R	500U
106-47-8 4-Chloroaniline U J	1000U	56-55-3 Benzo(a)anthracene	200U
87-68-3 Hexachlorobutadiene	300U	117-81-7 bis (2-Ethylhexyl) Phthalate J	360B
59-50-7 4-Chloro-3-Methylphenol	200U	218-01-9 Chrysene	200U
11-57-6 2-Methylnaphthalene	100U	117-84-0 Di-n-Octyl Phthalate	200U
77-47-4 Hexachlorocyclopentadiene	800U	205-99-2 Benzo(b)Fluoranthene	200U
88-06-2 2,4,6-Trichlorophenol	100U	207-08-9 Benzo(k)Fluoranthene U J	600U
15-95-4 2,4,5-Trichlorophenol	100U	50-32-8 Benzo(a)Pyrene	200U
11-58-7 2-Chloronaphthalene	200U	193-39-5 Indeno(1,2,3- α)Pyrene	200U
88-74-4 2-Nitroaniline - R	1000U	53-70-3 Dibenzo(a,h)Anthracene	200U
131-11-3 Dimethyl Phthalate	200U	100U	200U
106-96-8 Acenaphthylene ~ U J	1000U	191-24-2 Benzo(g,h,i)Perylene	200U
99-09-2 3-Nitroaniline - U J	1000U	(1) - Cannot be separated from diphenylamine	

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Sample Number

H1847

PC
10/17/87

Organics Analysis Data Sheet
(Page 3)

ANALYSIS NOT REQUIRED

Pesticide/PCBs

Concentration: Low Medium (Circle One)

Date Extracted/Prepared:

Date Analyzed:

Conc/Dil Factor:

CAS Number	ug/l or ug/Kg (Circle One)
319-84-6	Alpha-BHC
319-85-7	Beta-BHC
319-86-8	Delta-BHC
58-89-9	Gamma-BHC (Lindane)
76-44-9	Heptachlor
309-00-2	Aldrin
1024-57-3	Heptachlor Epoxide
959-98-8	Endosulfan I
60-57-1	Dieldrin
72-55-9	4, 4'-DDE
72-20-8	Endrin
33213-65-9	Endosulfan II
72-54-8	4, 4'-DDD
7421-93-4	Endrin Aldehyde
1031-07-8	Endosulfan Sulfate
50-29-3	4, 4'-DDT
72-43-5	Methoxychlor
53494-70-5	Endrin Ketone
57-74-9	Chlordane
8001-35-2	Toxaphene
12674-11-2	Aroclor-1016
11104-28-2	Aroclor-1221
11141-16-5	Aroclor-1232
53469-21-9	Aroclor-1242
12672-29-6	Aroclor-1248
11097-69-1	Aroclor-1254
11096-82-5	Aroclor-1260

V_i = Volume of extract injected (ul)

V_s = Volume of water extracted (ml)

W_s = Weight of sample extracted (g)

V_t = Volume of total extract (ul)

V_s _____ or W_s _____ V_t _____ V_i _____

3 290

Sample Number
H1847

2/2/84

Organics Analysis Data Sheet
(Page 4)

Tentatively Identified Compounds

No volatile compounds found.

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or ug/kg)
1. 96-48-0	2(3H)-Furanone, dihydro-	BNA	580	200JB
2. 628-68-2	Ethanol, 2,2'-oxybis-, diacetate	BNA	695	2000JB
3. --	Unknown fatty acid	BNA	1468	100J
4. --	Unknown amide	BNA	1709	600J
5. 103-23-1	Hexanedioic acid, bis(2-ethylhexyl)ester	BNA	1723	300JB
6. --	Unknown	BNA	1770	200J
7. --	Unknown	BNA	1864	400J
8. --	Unknown saturated aliphatic hydrocarbon	BNA	1867	400J
9. --	Unknown	BNA	1959	800J
10. --	Unknown saturated aliphatic hydrocarbon	BNA	1972	2000J
11. --	Unknown aldehyde	BNA	1977	2000J
12. --	Unknown saturated aliphatic hydrocarbon	BNA	2112	800J
13. --	Unknown	BNA	2121	1000J
14. --				
15. --				
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Sample Number
H1849

RE
6/1/85

Organics Analysis Data Sheet
(Page 1)

Laboratory Name: CE-EMSI

1 b Sample ID No: 050185V10 (VOA) 051885C09 (BNA)

Sample Matrix: Soil

Data Release Authorized By:

Richard Klontos

Case No: 4242

QC Report No: 4242

Contract No: 68-01-6783

Date Sample Received: 24-APR-85

Volatile Compounds

Concentration: Low
Date Extracted/Prepared: NA
Date Analyzed: 01-MAY-85
Conc/Dil Factor: NA pH 8.0
Percent Moisture: 65.84
Percent Moisture (Decanted): NA

S number		ug/l or ug/Kg (Circle one)	CAS Number		ug/l or ug/Kg (Circle one)
-87-3	Chloromethane	70U	79-34-5	1,1,2,2-Tetrachloroethane <input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>
-83-9	Bromomethane	6U	78-87-5	1,2-Dichloropropane <input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>
75-01-4	Vinyl Chloride <input checked="" type="checkbox"/>	70U	10061-02-6	Trans-1,3-Dichloropropene	<input checked="" type="checkbox"/>
--00-3	Chloroethane	6U	79-01-6	Trichloroethene	<input checked="" type="checkbox"/>
-09-2	Methylene Chloride <input checked="" type="checkbox"/>	72B	124-48-1	Dibromochloromethane	<input checked="" type="checkbox"/>
67-64-1	Acetone	37B	79-00-5	1,1,2-Trichloroethane	<input checked="" type="checkbox"/>
75-15-0	Carbon Disulfide	10U	71-43-2	Benzene	<input checked="" type="checkbox"/>
-35-4	1,1-Dichloroethene	9U	10061-01-5	cis-1,3-Dichloropropene	<input checked="" type="checkbox"/>
--34-3	1,1-Dichloroethane	8U	110-75-8	2-Chloroethylvinylether <input checked="" type="checkbox"/>	20U
156-60-5	Trans-1,2-Dichloroethene	6U	75-25-2	Bromoform	<input checked="" type="checkbox"/>
-66-3	Chloroform	3U	591-78-6	2-Hexanone <input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>
-7-06-2	1,2-Dichloroethane	3U	108-10-1	4-Methyl-2-Pentanone	<input checked="" type="checkbox"/>
78-93-3	2-Butanone <input checked="" type="checkbox"/>	3U	127-18-4	Tetrachloroethene	<input checked="" type="checkbox"/>
77-55-6	1,1,1-Trichloroethane <input checked="" type="checkbox"/>	9U	108-88-3	Toluene <input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>
-23-5	Carbon Tetrachloride	9U	108-90-7	Chlorobenzene	<input checked="" type="checkbox"/>
106-05-4	Vinyl Acetate <input checked="" type="checkbox"/>	6U	100-41-4	Ethylbenzene	<input checked="" type="checkbox"/>
75-27-4	Bromodichloromethane	3U	100-42-5	Styrene	<input checked="" type="checkbox"/>
				Total Xylenes	<input checked="" type="checkbox"/>

Data Reporting Qualifiers

For reporting results to EPA, the following results qualifiers are used.

J If the result is a value greater than or equal to the detection limit, report the value

C This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides $\geq 10 \text{ ng}/\text{ul}$ in the final extract should be confirmed by GC/MS.

U Compound was analyzed for but not detected. The number is the minimum attainable detection limit for the sample.

B This flag is used when the analyte is found in the blank as well as a sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action.

J Indicates an estimated value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicates the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero (e.g. 10U).

S Indicates compound was spiked into the sample for the matrix recovery study. The amounts that were spiked into and recovered from the sample are reported on Form III.

Environmental Protection Agency. CLP Sample Management Office.
P.O. Box 818. Alexandria, Virginia 22313 703/557-2490

Sample Number
H1849

2/2/85

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Organics Analysis Data Sheet
(Page 2)

Semivolatile Compounds

Concentration: Medium
Date Extracted/Prepared: 29-APR-85
Date Analyzed: 18-MAY-85
(Conc/Dil Factor: 5.0)

CAS Number	ug/l or ug/Kg (Circle one)	CAS Number	ug/l or ug/Kg (Circle one)		
62-75-9	N-Nitrosodimethylamine	100000U	83-32-9	Acenaphthene	50000U
108-95-2	Phenol - R	30000U	51-28-5	2,4-Dinitrophenol - R	300000U
1-53-3	Aniline R	60000U	100-02-7	4-Nitrophenol - R	400000U
1-44-4	bis(2-Chloroethyl)Ether R	60000U	132-64-9	Dibenzofuran	40000U
95-57-8	2-Chlorophenol	30000U	121-14-2	2,4-Dinitrotoluene	100000U
141-73-1	1,3-Dichlorobenzene	30000U	606-20-2	2,6-Dinitrotoluene	100000U
6-46-7	1,4-Dichlorobenzene	30000U	84-66-2	Diethylphthalate	50000U
100-51-6	Benzyl Alcohol	100000U	7005-72-3	4-Chlorophenyl-phenylether	20000U
95-50-1	1,2-Dichlorobenzene	30000U	86-73-7	Fluorene	30000U
48-7	2-Methylphenol	30000U	100-01-6	4-Nitroaniline - R	100000U
59-638-32-9	bis (2-Chloroisopropyl) Ether	30000U	534-52-1	4,6-Dinitro-2-Methylphenol	300000U
106-44-5	4-Methylphenol	200000U	86-30-6	N-Nitrosodiphenylamine (1)	100000U
1-64-7	N-Nitroso-Di-n-Propylamine	100000U	101-55-3	4-Ethoxyphenylphenylether - R	70000U
72-1	Hexachloroethane	40000U	118-74-1	Hexachlorobenzene - R	40000U
98-95-3	Nitrobenzene	100000U	87-66-5	Pentachlorophenol	40000U
70-59-1	Isophorone	60000U	85-01-8	Phenanthrene	20000U
1-75-5	2-Nitrophenol	40000U	120-12-7	Anthracene - R	30000U
105-67-9	2,4-Dimethylphenol	20000U	84-74-2	Di-n-Butylphthalate - R	200000U
65-85-0	Benzoic Acid R	100000U	206-44-0	Fluoranthene	50000U
1-91-1	bis (2-Chloroethoxy) Methane	70000U	92-87-5	Benzidine - R	200000U
90-63-2	2,4-Dichlorophenol	20000U	129-00-0	Pyrene	100000U
120-82-1	1,2,4-Trichlorobenzene	50000U	85-68-7	Butylbenzylphthalate	60000U
7-20-3	Naphthalene	50000U	91-94-1	3,3'-Dichlorobenzidine - R	100000U
16-47-8	4-Chloroaniline R	300000U	56-55-3	Benzo(a)anthracene	30000U
87-68-3	Hexachlorobutadiene	60000U	117-81-7	bis (2-Ethylhexyl) Phthalate R	30000U
90-50-7	4-Chloro-3-Methylphenol	40000U	218-01-9	Chrysene	30000U
57-6	2-Methylnaphthalene	20000U	117-84-0	Di-n-octyl Phthalate	30000U
77-47-4	Hexachlorocyclopentadiene	200000U	205-99-2	Benzo(b)Fluoranthene	100000U
88-06-2	2,4,6-Trichlorophenol	20000U	207-08-9	Benzo(k)Fluoranthene	40000U
195-4	2,4,5-Trichlorophenol	20000U	50-32-8	Benzo(a)Pyrene	30000U
1-58-7	2-Chlorosanthalene	40000U	193-39-5	Indeno(1,2,3-cd)Pyrene	30000U
88-74-4	2-Nitroaniline - R	200000U	53-70-3	Dibenzo(a,h)Anthracene	30000U
71-11-3	Dimethyl Phthalate	40000U	191-24-2	Benzo(g,h,i)Perylene	40000U
19-96-8	Acenaphthylene R	20000U			
95-09-2	3-Nitroaniline R	200000U			

(1)- Cannot be separated from diphenylamine

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Organics Analysis Data Sheet
(Page 3)

ANALYSIS NOT REQUIRED

Pesticide/PCBs

Concentration: Low Medium (Circle One)

Date Extracted/Prepared: _____

Date Analyzed: _____

Conc/Dil Factor: _____

CAS Number		ug/l or ug/Kg (Circle One)
319-84-6	Alpha-BHC	
319-85-7	Beta-BHC	
319-86-8	Delta-BHC	
58-89-9	Gamma-BHC (Lindane)	
76-44-8	Heptachlor	
309-00-2	Aldrin	
1024-57-3	Heptachlor Epoxide	
959-98-8	Endosulfan I	
60-57-1	Dieldrin	
72-55-9	4, 4'-DDE	
72-20-8	Endrin	
33213-65-9	Endosulfan II	
72-54-8	4, 4'-DDD	
7421-93-4	Endrin Aldehyde	
1031-07-8	Endosulfan Sulfate	
50-29-3	4, 4'-DDT	
72-43-5	Methoxychlor	
53494-70-5	Endrin Ketone	
57-74-9	Chlordane	
8001-35-2	Taxaphene	
12674-11-2	Aroclor-1016	
11104-28-2	Aroclor-1221	
11141-16-5	Aroclor-1232	
53469-21-9	Aroclor-1242	
12672-29-6	Aroclor-1248	
11097-69-1	Aroclor-1254	
11096-82-5	Aroclor-1260	

V_i = Volume of extract injected (ul)

V_s = Volume of water extracted (ml)

W_s = Weight of sample extracted (g)

V_t = Volume of total extract (ul)

V_s _____ or W_s _____ V_t _____ V_i _____

Sample Number
H1849

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9/6/85

Organics Analysis Data Sheet
(Page 4)

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or ug/kg)
1. --	x,y,z-Trimethylcyclohexane	VOA	380	9J
2. --	C10 H18 Cyclic hydrocarbon	VOA	419	8J
3. --	C10 H18 Cyclic hydrocarbon & Unknown	VOA	445	10J
4. --	Trap artifact	VOA	457	
5. --	C10 H18 Cyclic hydrocarbon	VOA	490	10J
6. --	Unknown	VOA	516	10J
7. --	Unknown	VOA	525	10J
8. --	C10 H20 Cyclic hydrocarbon	VOA	554	60J
9. 54676-39-0	Cyclohexane, 2-butyl-1,1,3-trimethyl-	BNA	941	8000J
10. --	Unknown saturated aliphatic hydrocarbon	BNA	974	8000J
11. --	C14 H28 Hydrocarbon	BNA	993	6000J
12. --	Unknown saturated aliphatic hydrocarbon	BNA	1015	5000J
13. --	Unknown saturated aliphatic hydrocarbon	BNA	1066	30000J
14. --	C15 H30 Hydrocarbon	BNA	1088	20000J
15. --	Unknown saturated aliphatic hydrocarbon	BNA	1137	30000J
16. --	Unknown saturated aliphatic hydrocarbon	BNA	1153	10000J
17. --	Unknown hydrocarbon (MW-224)	BNA	1161	8000J
18. --	Unknown saturated aliphatic hydrocarbon	BNA	1219	8000J
19. --	Unknown saturated aliphatic hydrocarbon	BNA	1281	5000J
20. --				
21. --				
22. --				
23. --				
24. --				
25. --				
26. --				
27. --				
28. --				
29. --				
30. --				

APPENDIX B
FINAL SITE INSPECTION FORM



POTENTIAL HAZARDOUS WASTE SITE
SITE INSPECTION REPORT
PART 1 - SITE LOCATION AND INSPECTION INFORMATION

01 STATE WY	02 SITE NUMBER D980807762
----------------	------------------------------

II. SITE NAME AND LOCATION

01 SITE NAME (Legal, common or descriptive name of site) Empire State Oil Company	02 STREET, ROUTE NO., OR SPECIFIC LOCATION IDENTIFIER 242 Amoretti Street				
03 CITY Thermopolis	04 STATE WY	05 ZIP CODE 82443	06 COUNTY Hot Springs	07 COUNTY CODE 017	08 CONG DIST 00
09 COORDINATES LATITUDE 43 38 21.	LONGITUDE 108 12 25.	10 TYPE OF OWNERSHIP (Check one) <input checked="" type="checkbox"/> A. PRIVATE <input type="checkbox"/> B. FEDERAL <input type="checkbox"/> C. STATE <input type="checkbox"/> D. COUNTY <input type="checkbox"/> E. MUNICIPAL <input type="checkbox"/> F. OTHER <input type="checkbox"/> G. UNKNOWN			

III. INSPECTION INFORMATION

01 DATE OF INSPECTION 6 / 25, 84	02 SITE STATUS <input type="checkbox"/> ACTIVE <input checked="" type="checkbox"/> INACTIVE	03 YEARS OF OPERATION 1938 - 1969	BEGINNING YEAR	ENDING YEAR	UNKNOWN
--	--	---	----------------	-------------	---------

04 AGENCY PERFORMING INSPECTION (Check all that apply)	(E&E)				
<input type="checkbox"/> A. EPA <input checked="" type="checkbox"/> B. EPA CONTRACTOR <input type="checkbox"/> C. MUNICIPAL <input type="checkbox"/> D. MUNICIPAL CONTRACTOR	(Name of firm)				(Name of firm)
<input type="checkbox"/> E. STATE <input type="checkbox"/> F. STATE CONTRACTOR <input type="checkbox"/> G. OTHER	(Name of firm)				(Name of firm)

05 CHIEF INSPECTOR Mark Mullis	06 TITLE FIT Environmental Scientist	07 ORGANIZATION E&E	08 TELEPHONE NO (303) 757-4984
09 OTHER INSPECTORS Ben Genes	10 TITLE FIT Environmental Engineer	11 ORGANIZATION E&E	12 TELEPHONE NO. (303) 757-4984
Teresa Sherman	FIT Environmental Engineer	E&E	(303) 757-4984
Chris Renda	FIT Hydrogeologist	E&E	(303) 757-4984
Susan Kennedy	FIT Environmental Scientist	E&E	(303) 757-4984
Ken Moll	FIT Geologist	E&E	(303) 757-4984
Steve Yarbrough	FIT Environmental Scientist	E&E	(303) 757-4984
13 SITE REPRESENTATIVES INTERVIEWED Doug Wornock	14 TITLE Administrator	15 ADDRESS Gottzsche Foundation P.O. Box 780	16 TELEPHONE NO (307) 864-3532
		Thermopolis, WY 82443	()
			()

17 ACCESS GAINED BY (Check one) <input checked="" type="checkbox"/> PERMISSION <input type="checkbox"/> WARRANT	18 TIME OF INSPECTION 0830	19 WEATHER CONDITIONS mostly clear skies, 70°F, light breeze
--	--------------------------------------	--

IV. INFORMATION AVAILABLE FROM

01 CONTACT Doug Wornock	02 OF (Agency/Organization) Gottzsche Foundation	03 TELEPHONE NO. (307) 864-3532		
04 PERSON RESPONSIBLE FOR SITE INSPECTION FORM Mark Mullis	05 AGENCY EPA Contractor	06 ORGANIZATION Ecology & Environment	07 TELEPHONE NO. (303) 757-4984	08 DATE 7 / 12 / 84
				MONTH DAY YEAR



**POTENTIAL HAZARDOUS WASTE SITE
SITE INSPECTION REPORT
PART 2 - WASTE INFORMATION**

IDENTIFICATION	
01 STATE	02 SITE NUMBER
WY	D980807762

II. WASTE STATES, QUANTITIES, AND CHARACTERISTICS

D1 PHYSICAL STATES (Check all that apply)		D2 WASTE QUANTITY AT SITE <small>(Measures of waste quantity; must be independent)</small>		D3 WASTE CHARACTERISTICS (Check all that apply)					
<input type="checkbox"/> A SOLID	<input type="checkbox"/> D E SLURRY	TONS _____		<input type="checkbox"/> A TOXIC	<input type="checkbox"/> E SOLUBLE	<input checked="" type="checkbox"/> I HIGHLY VOLATILE			
<input type="checkbox"/> B POWDER, FINES	<input type="checkbox"/> E F LIQUID	CUBIC YARDS _____		<input type="checkbox"/> B CORROSIVE	<input type="checkbox"/> F INFECTIOUS	<input type="checkbox"/> J EXPLOSIVE			
<input type="checkbox"/> C SLUDGE	<input type="checkbox"/> G GAS	NO. OF DRUMS _____		<input type="checkbox"/> C RADIOACTIVE	<input type="checkbox"/> G FLAMMABLE	<input type="checkbox"/> K REACTIVE			
<input type="checkbox"/> D OTHER _____ <small>(Specify)</small>				<input type="checkbox"/> D PERSISTENT	<input type="checkbox"/> H IGNITABLE	<input type="checkbox"/> L INCOMPATIBLE			
						<input type="checkbox"/> M NOT APPLICABLE			

III. WASTE TYPE

CATEGORY	SUBSTANCE NAME	Q1 GROSS AMOUNT	Q2 UNIT OF MEASURE	Q3 COMMENTS
SLU	SLUDGE			
OLW	OILY WASTE	unknown		waste spilled on ground and disposed of in pits
SOL	SOLVENTS			
PSD	PESTICIDES			
OCC	OTHER ORGANIC CHEMICALS			
IOC	INORGANIC CHEMICALS			
ACD	ACIDS			
BAS	BASES			
MES	HEAVY METALS			

IV. HAZARDOUS SUBSTANCES (See Appendix for most frequently cited CAS Numbers)

V. FEEDSTOCKS (See Appendix B: Gas Blends)

CATEGORY	O1 FEEDSTOCK NAME	O2 CAS NUMBER	CATEGORY	O1 FEEDSTOCK NAME	O2 CAS NUMBER
FDS			FDS		
FDS	None		FDS		
FDS			FDS		
FDS			FDS		

VL SOURCES OF INFORMATION (the specific references e.g., state the sample analysis reports)

Preliminary Assessment, 5/18/84, State of Wyoming, Dept. of Environmental Quality,
Solid Waste, TDD R8-8405-31
Site Visit 6/25/84, Ecology and Environment, Inc., TDD R8-8412-07
Sampling Activities Report 4/20/85, Ecology and Environment, Inc., TDD R8-8503-20

POTENTIAL HAZARDOUS WASTE SITE SITE INSPECTION REPORT		I. IDENTIFICATION
		01 STATE 02 SITE NUMBER WY D980807762
PART 3 - DESCRIPTION OF HAZARDOUS CONDITIONS AND INCIDENTS		
II. HAZARDOUS CONDITIONS AND INCIDENTS		
<p>01 <input type="checkbox"/> A GROUNDWATER CONTAMINATION APPROX. 02 <input type="checkbox"/> OBSERVED (DATE _____) 03 POPULATION POTENTIALLY AFFECTED <u>5000</u> 04 NARRATIVE DESCRIPTION <input type="checkbox"/> POTENTIAL <input type="checkbox"/> ALLEGED</p> <p>Spills which may have occurred during the life time of the refinery and possible leakage from disposal pits may have contaminated the underlying alluvial aquifer. See Section IV under Part 3.</p>		
<p>01 <input type="checkbox"/> B SURFACE WATER CONTAMINATION APPROX. 02 <input type="checkbox"/> OBSERVED (DATE <u>4/21/85</u>) 03 POPULATION POTENTIALLY AFFECTED <u>5000</u> 04 NARRATIVE DESCRIPTION <input type="checkbox"/> POTENTIAL <input checked="" type="checkbox"/> ALLEGED</p> <p>Toluene detected at 57 ug/l in Bighorn River. Mercury detected at 1.7 ug/l in pond on site. Aluminum and manganese detected in pond. These concentrations are above the concentration in the background sample.</p>		
<p>01 <input type="checkbox"/> C CONTAMINATION OF AIR 02 <input type="checkbox"/> OBSERVED (DATE _____) 03 POPULATION POTENTIALLY AFFECTED _____ 04 NARRATIVE DESCRIPTION <input type="checkbox"/> POTENTIAL <input type="checkbox"/> ALLEGED</p> <p>None known</p>		
<p>01 <input type="checkbox"/> D FIRE/EXPLOSIVE CONDITIONS 02 <input type="checkbox"/> OBSERVED (DATE _____) 03 POPULATION POTENTIALLY AFFECTED _____ 04 NARRATIVE DESCRIPTION <input type="checkbox"/> POTENTIAL <input type="checkbox"/> ALLEGED</p> <p>None known</p>		
<p>01 <input type="checkbox"/> E DIRECT CONTACT 02 <input type="checkbox"/> OBSERVED (DATE _____) 03 POPULATION POTENTIALLY AFFECTED <u>5000</u> with <u>in 1 mile</u> 04 NARRATIVE DESCRIPTION <input type="checkbox"/> POTENTIAL <input type="checkbox"/> ALLEGED</p> <p>Petroleum derivatives, including asphalt, lie on the surface of the ground in certain areas. Site is accessible to humans and animals. See Section IV under Part 3.</p>		
<p>01 <input type="checkbox"/> F CONTAMINATION OF SOIL 02 <input type="checkbox"/> OBSERVED (DATE <u>6/25/84</u>) 03 AREA POTENTIALLY AFFECTED <u>40</u> (Acres) 04 NARRATIVE DESCRIPTION <input type="checkbox"/> POTENTIAL <input type="checkbox"/> ALLEGED</p> <p>Spills from above ground tanks and in refining area were evident during FIT visit. During 31 years of operation, the possibility existed for significant soil contamination. The soil sample collected near the refining area contained elevated cadmium, lead, ethylbenzene, methylene chloride, xylenes, and di-n-butylphthalate.*</p>		
<p>01 <input type="checkbox"/> G DRINKING WATER CONTAMINATION APPROX. 02 <input type="checkbox"/> OBSERVED (DATE _____) 03 POPULATION POTENTIALLY AFFECTED <u>5000</u> 04 NARRATIVE DESCRIPTION <input type="checkbox"/> POTENTIAL <input type="checkbox"/> ALLEGED</p> <p>Well field from which Thermopolis and vicinity draws much of its drinking water is 1600 feet WSW of the refinery property across the Big Horn River. The wells are only 30 feet deep. See Section IV under Part 3.</p>		
<p>01 <input type="checkbox"/> H WORKER EXPOSURE/INJURY 02 <input type="checkbox"/> OBSERVED (DATE _____) 03 WORKERS POTENTIALLY AFFECTED _____ 04 NARRATIVE DESCRIPTION <input type="checkbox"/> POTENTIAL <input type="checkbox"/> ALLEGED</p> <p>None known</p>		
<p>01 <input type="checkbox"/> I. POPULATION EXPOSURE/INJURY 02 <input type="checkbox"/> OBSERVED (DATE _____) 03 POPULATION POTENTIALLY AFFECTED _____ 04 NARRATIVE DESCRIPTION <input type="checkbox"/> POTENTIAL <input type="checkbox"/> ALLEGED</p> <p>None known</p>		



POTENTIAL HAZARDOUS WASTE SITE
SITE INSPECTION REPORT

PART 3 - DESCRIPTION OF HAZARDOUS CONDITIONS AND INCIDENTS

I. IDENTIFICATION

01 STATE WY	02 SITE NUMBER D980807762
----------------	------------------------------

II. HAZARDOUS CONDITIONS AND INCIDENTS (Continued)

01 J. DAMAGE TO FLORA
04 NARRATIVE DESCRIPTION

02 OBSERVED (DATE 6/25/84) POTENTIAL ALLEGED

Indication of vegetation stress observed by FIT. Vegetation is largely absent from soil now covering former disposal pits.

01 K. DAMAGE TO FAUNA Dead fish
04 NARRATIVE DESCRIPTION (include name(s) of species)

02 OBSERVED (DATE 3/30/84) POTENTIAL ALLEGED

Observed by WDEQ

Not observed by FIT on 6/25/84

01 L. CONTAMINATION OF FOOD CHAIN
04 NARRATIVE DESCRIPTION

02 OBSERVED (DATE _____) POTENTIAL ALLEGED

None known

01 M. UNSTABLE CONTAINMENT OF WASTES
(Soil Runoff Standing toward leaking drums)
02 OBSERVED (DATE _____) POTENTIAL ALLEGED
03 POPULATION POTENTIALLY AFFECTED Approx. 5000
04 NARRATIVE DESCRIPTION

Site dismantled and unlined pits covered with soil. See Section IV.

01 N. DAMAGE TO OFFSITE PROPERTY
04 NARRATIVE DESCRIPTION

02 OBSERVED (DATE _____) POTENTIAL ALLEGED

None known

01 O. CONTAMINATION OF SEWERS STORM DRAINS WWTPs
04 NARRATIVE DESCRIPTION

02 OBSERVED (DATE _____) POTENTIAL ALLEGED

None known

01 P. ILLEGAL/UNAUTHORIZED DUMPING
04 NARRATIVE DESCRIPTION

02 OBSERVED (DATE _____) POTENTIAL ALLEGED

None known

05 DESCRIPTION OF ANY OTHER KNOWN, POTENTIAL, OR ALLEGED HAZARDS

None

II. TOTAL POPULATION POTENTIALLY AFFECTED: Approx. 5000

IV. COMMENTS

For A,B,E,G, and M, above, the population potentially affected is shown, but the location of the wells and surface water intake make the possibility for a significant health hazard moderate.

V. SOURCES OF INFORMATION (Cite specific references, e.g., State Baseline analysis reports.)

Preliminary Assessment, 5/18/84, State of Wyoming, Department of Environmental Quality, Division of Solid Waste, TDD R8-8405-31
Site Visit, 6/25/84, Ecology and Environment, Inc.
Report of Sampling Activities 4/20/85, TDD R8-8503-20



POTENTIAL HAZARDOUS WASTE SITE
SITE INSPECTION
PART 4 - PERMIT AND DESCRIPTIVE INFORMATION

I. IDENTIFICATION
01 STATE | 02 SITE NUMBER
WY | D980807762

II. PERMIT INFORMATION

01 TYPE OF PERMIT ISSUED (Check all that apply)	02 PERMIT NUMBER	03 DATE ISSUED	04 EXPIRATION DATE	05 COMMENTS
<input type="checkbox"/> A. NPDES				
<input type="checkbox"/> B. UIC				
<input type="checkbox"/> C. AIR				
<input type="checkbox"/> D. RCRA				
<input type="checkbox"/> E. RCRA INTERIM STATUS				
<input type="checkbox"/> F. SPCC PLAN				
<input type="checkbox"/> G. STATE (Specify)				
<input type="checkbox"/> H. LOCAL (Specify)				
<input type="checkbox"/> I. OTHER (Specify)				
<input checked="" type="checkbox"/> J. NONE				

III. SITE DESCRIPTION

01 STORAGE/DISPOSAL (Check all that apply)	02 AMOUNT	03 UNIT OF MEASURE	04 TREATMENT (Check all that apply)	05 OTHER
<input checked="" type="checkbox"/> A. SURFACE IMPOUNDMENT	unknown		<input type="checkbox"/> A. INCINERATION	<input checked="" type="checkbox"/> A. BUILDINGS ON SITE
<input type="checkbox"/> B. PILES			<input type="checkbox"/> B. UNDERGROUND INJECTION	
<input type="checkbox"/> C. DRUMS, ABOVE GROUND			<input type="checkbox"/> C. CHEMICAL/PHYSICAL	
<input checked="" type="checkbox"/> D. TANK, ABOVE GROUND			<input type="checkbox"/> D. BIOLOGICAL	
<input type="checkbox"/> E. TANK, BELOW GROUND			<input type="checkbox"/> E. WASTE OIL PROCESSING	
<input type="checkbox"/> F. LANDFILL			<input type="checkbox"/> F. SOLVENT RECOVERY	
<input type="checkbox"/> G. LANDFARM			<input type="checkbox"/> G. OTHER RECYCLING/RECOVERY	
<input type="checkbox"/> H. OPEN DUMP			<input type="checkbox"/> H. OTHER (Specify)	
<input type="checkbox"/> I. OTHER (Specify)				

07 COMMENTS

The facility is no longer active. Tanks have been removed and waste pits covered.

IV. CONTAINMENT

01 CONTAINMENT OF WASTES (Check one)	02 COMMENTS	03 CONTAINMENT OF WASTES (Check one)	04 COMMENTS
<input type="checkbox"/> A. ADEQUATE, SECURE		<input checked="" type="checkbox"/> B. MODERATE	<input checked="" type="checkbox"/> C. INADEQUATE, POOR

02 DESCRIPTION OF DRUMS, DIKING, LINERS, BARRIERS, ETC.

Disposal pits contain no liners, leakage observed around removed above ground tanks

V. ACCESSIBILITY

01 WASTE EASILY ACCESSIBLE YES NO

02 COMMENTS

Fence about perimeter, but gate was unlocked and unattended.

VI. SOURCES OF INFORMATION (Give specific references, e.g. state files, sample analysis reports)

Preliminary Assessment, 5/18/84, State of Wyoming, Dept. of Environmental Quality.
Site Visit, 6/25/84, FIT, Ecology and Environment, Inc.



POTENTIAL HAZARDOUS WASTE SITE
SITE INSPECTION REPORT
PART 5 - WATER, DEMOGRAPHIC, AND ENVIRONMENTAL DATA

I IDENTIFICATION
01 STATE: WY 02 SITE NUMBER: D980807762

A DRINKING WATER SUPPLY

TYPE OF DRINKING SUPPLY <small>(Check one applicable)</small>		02 STATUS	For hardness & TDS MONITORED		03 DISTANCE TO SITE
COMMUNITY	SURFACE A. <input checked="" type="checkbox"/> B. <input type="checkbox"/> NON-COMMUNITY C. <input type="checkbox"/> D. <input checked="" type="checkbox"/>	WELL B. <input checked="" type="checkbox"/> D. <input type="checkbox"/>	ENDANGERED A. <input type="checkbox"/> B. <input type="checkbox"/> D. <input type="checkbox"/>	AFFECTED E. <input type="checkbox"/> F. <input type="checkbox"/>	A 0.2 (mi) B _____ (mi)

B GROUNDWATER

GROUNDWATER USE IN VICINITY (Check one)		03 COMMERCIAL INDUSTRIAL IRRIGATION <small>(Limited drinking sources available)</small>		04 NOT USED, UNUSEABLE	
<input checked="" type="checkbox"/> A ONLY SOURCE FOR DRINKING	<input checked="" type="checkbox"/> B DRINKING <small>(Other sources available)</small> COMMERCIAL INDUSTRIAL, IRRIGATION <small>(No other water sources available)</small>				

POPULATION SERVED BY GROUND WATER		05 DIRECTION OF GROUNDWATER FLOW	03 DISTANCE TO NEAREST DRINKING WATER WELL	0.2 (mi)
DEPTH TO GROUNDWATER	15 (ft)	SE toward river	06 DEPTH TO AQUIFER OF CONCERN	07 POTENTIAL YIELD OF AQUIFER
			13 (ft)	1.4 million (gpd) <input checked="" type="checkbox"/> NO <input type="checkbox"/>

DESCRIPTION OF WELLS (including usage, depth, and location relative to population and buildings)

3 municipal wells, 30 feet deep, supply 1000 gpm to city system. The wells are on the floodplain across the river and directly south of the town of Thermopolis.

DISCHARGE AREA		11 DISCHARGE AREA	
YES	COMMENTS	YES	COMMENTS
<input checked="" type="checkbox"/>	In spring, river recharges alluvial aquifer	<input type="checkbox"/>	In fall, aquifer recharges river.

C SURFACE WATER

SURFACE WATER USE (Check one)		12 AFFECTED		DISTANCE TO SITE	
<input checked="" type="checkbox"/> A RESERVOIR, RECREATION DRINKING WATER SOURCE	<input type="checkbox"/> B IRRIGATION, ECONOMICALLY IMPORTANT RESOURCES	<input type="checkbox"/> C COMMERCIAL INDUSTRIAL	<input type="checkbox"/> D NOT CURRENTLY USED		

D AFFECTED/POTENTIALLY AFFECTED BODIES OF WATER

NAME	AFFECTED	DISTANCE TO SITE
Big Horn River	<input type="checkbox"/>	0 (mi)
	<input type="checkbox"/>	(mi)
	<input type="checkbox"/>	(mi)

E DEMOGRAPHIC AND PROPERTY INFORMATION

TOTAL POPULATION WITHIN			02 DISTANCE TO NEAREST POPULATION
A 1 MILE OF SITE 4500 NO OF PERSONS	B TWO (2) MILES OF SITE 5000 NO OF PERSONS	C THREE (3) MILES OF SITE 5000 NO OF PERSONS	Across the street (mi)

NUMBER OF BUILDINGS WITHIN TWO (2) MILES OF SITE	04 DISTANCE TO NEAREST OFF-SITE BUILDING
Entire town of Thermopolis	Across the street (mi)

F POPULATION WITHIN VICINITY OF SITE (Provide narrative description of nature of population within vicinity of site e.g. rural, village, densely populated urban area)

Approximately 5000 people live in the vicinity of Thermopolis.

 POTENTIAL HAZARDOUS WASTE SITE SITE INSPECTION REPORT PART 5 - WATER, DEMOGRAPHIC, AND ENVIRONMENTAL DATA		I. IDENTIFICATION D1 STATE WY D2 SITE NUMBER D980807762	
VI. ENVIRONMENTAL INFORMATION			
D1 PERMEABILITY OF UNSATURATED ZONE (Check one)			
<input type="checkbox"/> A. $10^{-6} - 10^{-8}$ cm/sec <input type="checkbox"/> B. $10^{-4} - 10^{-6}$ cm/sec <input checked="" type="checkbox"/> C. $10^{-4} - 10^{-3}$ cm/sec <input type="checkbox"/> D. GREATER THAN 10^{-3} cm/sec			
D2 PERMEABILITY OF BEDROCK (Check one)			
<input type="checkbox"/> A. IMPERMEABLE ($\text{flow} < 10^{-6}$ cm/sec) <input checked="" type="checkbox"/> B. RELATIVELY IMPERMEABLE ($10^{-4} - 10^{-6}$ cm/sec) <input type="checkbox"/> C. RELATIVELY PERMEABLE ($10^{-2} - 10^{-4}$ cm/sec) <input type="checkbox"/> D. VERY PERMEABLE ($\text{flow} > 10^{-3}$ cm/sec)			
D3 DEPTH TO BEDROCK <u>Approx. 40</u> (ft)	D4 DEPTH OF CONTAMINATED SOIL ZONE <u>unknown</u> (ft)	D5 SOIL pH <u>unknown</u>	
D6 NET PRECIPITATION <u>-31</u> (in)	D7 ONE YEAR 24 HOUR RAINFALL <u>Approx. 1.25</u> (in)	D8 SLOPE SITE SLOPE <u><1</u> %	DIRECTION OF SITE SLOPE <u>SE</u> TERRAIN AVERAGE SLOPE <u>1</u> %
D9 FLOOD POTENTIAL Dam immed. upstream SITE IS IN _____ YEAR FLOODPLAIN	D10 SITE IS ON BARRIER ISLAND, COASTAL HIGH HAZARD AREA, RIVERINE FLOODWAY		
D11 DISTANCE TO WETLANDS (3 acre minimum): ESTUARINE <u>A. N/A</u> (mi)	OTHER <u>B.</u> (mi)	D12 DISTANCE TO CRITICAL HABITAT (of endangered species): ENDANGERED SPECIES _____	
D13 LAND USE IN VICINITY DISTANCE TO: COMMERCIAL/INDUSTRIAL <u>A. 1/4</u> (mi)			
RESIDENTIAL AREAS, NATIONAL/STATE PARKS, FORESTS, OR WILDLIFE RESERVES			
AGRICULTURAL LANDS PRIME AG LAND AG LAND			
B <u>1</u> (mi) C <u>unknown</u> (mi) D <u>unknown</u> (mi)			
D14 DESCRIPTION OF SITE IN RELATION TO SURROUNDING TOPOGRAPHY Refinery is on the Bighorn River floodplain. Site has an average slope of 1% to the southeast.			
VII. SOURCES OF INFORMATION (Cite specific references, e.g., state files, sample analyses, reports)			
U.S.G.S. supplied some information on the geology of the Thermopolis area (307) 772-2716. U.S.G.S. 7.5' topographic map, Thermopolis Quad, 1:24,000 (307)772-2716 Steve Krysiak, Water and Sewer Superintendent, Town of Thermopolis (307) 864-2658			



POTENTIAL HAZARDOUS WASTE SITE
SITE INSPECTION REPORT
PART 6 - SAMPLE AND FIELD INFORMATION

I. IDENTIFICATION	
O1 STATE	O2 SITE NUMBER
WY	D980807762

II. SAMPLES TAKEN

SAMPLE TYPE	O1 NUMBER OF SAMPLES TAKEN	O2 SAMPLES SENT TO	O3 ESTIMATED DATE RESULTS AVAILABLE
GROUNDWATER			
SURFACE WATER	3	Organic - Combustion Eng., Ventura Cty, CA Inorganic - Chemtech Cons. Group LTD., NY, NY	
WASTE			
AIR			
RUNOFF			
SPILL			
SOIL	4	Organic - Combustion Eng., Ventura Cty, CA Inorganic - Chemtech Cons. Group LTD., NY, NY	
VEGETATION		Met Hazard - Rocky Mtn Analytical Lab, Denver, CO	
OTHER Sediment	3	Organic - Combustion Eng., Ventura Cty, CA Inorganic - Chemtech Cons. Group LTD., NY, NY	

III. FIELD MEASUREMENTS TAKEN

O1 TYPE	O2 COMMENTS
Organic Vapor Detection (HNU)	Downgradient surface water 0.8 ppm above background Downgradient sediment 0.8 ppm above background Disposal pit 0.4 - 0.8 ppm above background Refining area 4.0 ppm above background

IV. PHOTOGRAPHS AND MAPS

O1 TYPE	O2 IN CUSTODY OF
<input checked="" type="checkbox"/> GROUND <input type="checkbox"/> AERIAL	Ecology and Environment, Inc. <small>(Name of organization or individual)</small>
O3 MAPS	O4 LOCATION OF MAPS
<input checked="" type="checkbox"/> YES <input type="checkbox"/> NO	Ecology and Environment, Inc.

V. OTHER FIELD DATA COLLECTED (Provide narrative description)

Temperature, pH, and specific conductance was taken of all surface water samples.
 Temperature was between 6 and 10°C
 pH was between 5.31 and 9.33
 Conductivity was between 4 and 1800 umhos/cm

VI. SOURCES OF INFORMATION (Cite specific references e.g. state files, sample analysis reports)

Sampling effort of 4/20/85



**POTENTIAL HAZARDOUS WASTE SITE
SITE INSPECTION REPORT
PART 7 - OWNER INFORMATION**

D1 STATE	D2 SITE NUMBER
WY	D980807762

II. CURRENT OWNER(S)

D1 NAME	D2 D+B NUMBER	D3 STREET ADDRESS (P O Box, RFD #, etc.)	D4 SIC CODE	D5 CITY	D6 STATE	D7 ZIP CODE	D8 NAME	D9 D+B NUMBER	D10 STREET ADDRESS (P O Box, RFD #, etc.)	D11 SIC CODE
Gottzsche Foundation		P.O. Box 790		Thermopolis	WY	82443				
D1 NAME	D2 D+B NUMBER	D3 STREET ADDRESS (P O Box, RFD #, etc.)	D4 SIC CODE	D5 CITY	D6 STATE	D7 ZIP CODE	D8 NAME	D9 D+B NUMBER	D10 STREET ADDRESS (P O Box, RFD #, etc.)	D11 SIC CODE
D1 NAME	D2 D+B NUMBER	D3 STREET ADDRESS (P O Box, RFD #, etc.)	D4 SIC CODE	D5 CITY	D6 STATE	D7 ZIP CODE	D8 NAME	D9 D+B NUMBER	D10 STREET ADDRESS (P O Box, RFD #, etc.)	D11 SIC CODE
D1 NAME	D2 D+B NUMBER	D3 STREET ADDRESS (P O Box, RFD #, etc.)	D4 SIC CODE	D5 CITY	D6 STATE	D7 ZIP CODE	D8 NAME	D9 D+B NUMBER	D10 STREET ADDRESS (P O Box, RFD #, etc.)	D11 SIC CODE
D1 NAME	D2 D+B NUMBER	D3 STREET ADDRESS (P O Box, RFD #, etc.)	D4 SIC CODE	D5 CITY	D6 STATE	D7 ZIP CODE	D8 NAME	D9 D+B NUMBER	D10 STREET ADDRESS (P O Box, RFD #, etc.)	D11 SIC CODE
D1 NAME	D2 D+B NUMBER	D3 STREET ADDRESS (P O Box, RFD #, etc.)	D4 SIC CODE	D5 CITY	D6 STATE	D7 ZIP CODE	D8 NAME	D9 D+B NUMBER	D10 STREET ADDRESS (P O Box, RFD #, etc.)	D11 SIC CODE
D1 NAME	D2 D+B NUMBER	D3 STREET ADDRESS (P O Box, RFD #, etc.)	D4 SIC CODE	D5 CITY	D6 STATE	D7 ZIP CODE	D8 NAME	D9 D+B NUMBER	D10 STREET ADDRESS (P O Box, RFD #, etc.)	D11 SIC CODE
D1 NAME	D2 D+B NUMBER	D3 STREET ADDRESS (P O Box, RFD #, etc.)	D4 SIC CODE	D5 CITY	D6 STATE	D7 ZIP CODE	D8 NAME	D9 D+B NUMBER	D10 STREET ADDRESS (P O Box, RFD #, etc.)	D11 SIC CODE
D1 NAME	D2 D+B NUMBER	D3 STREET ADDRESS (P O Box, RFD #, etc.)	D4 SIC CODE	D5 CITY	D6 STATE	D7 ZIP CODE	D8 NAME	D9 D+B NUMBER	D10 STREET ADDRESS (P O Box, RFD #, etc.)	D11 SIC CODE

III. PREVIOUS OWNER(S) (list most recent first)

D1 NAME	D2 D+B NUMBER	D3 STREET ADDRESS (P O Box, RFD #, etc.)	D4 SIC CODE	D5 CITY	D6 STATE	D7 ZIP CODE	D8 NAME	D9 D+B NUMBER
Ashland Oil Company		Ashland Drive	29111					
D1 NAME	D2 D+B NUMBER	D3 STREET ADDRESS (P O Box, RFD #, etc.)	D4 SIC CODE	D5 CITY	D6 STATE	D7 ZIP CODE	D8 NAME	D9 D+B NUMBER
D1 NAME	D2 D+B NUMBER	D3 STREET ADDRESS (P O Box, RFD #, etc.)	D4 SIC CODE	D5 CITY	D6 STATE	D7 ZIP CODE	D8 NAME	D9 D+B NUMBER
		unknown (out of business)						
D1 NAME	D2 D+B NUMBER	D3 STREET ADDRESS (P O Box, RFD #, etc.)	D4 SIC CODE	D5 CITY	D6 STATE	D7 ZIP CODE	D8 NAME	D9 D+B NUMBER
D1 NAME	D2 D+B NUMBER	D3 STREET ADDRESS (P O Box, RFD #, etc.)	D4 SIC CODE	D5 CITY	D6 STATE	D7 ZIP CODE	D8 NAME	D9 D+B NUMBER

V. SOURCES OF INFORMATION (Check specific references e.g. state file sample analysis reports)

Preliminary Assessment, completed 5/18/84 by Wyoming Department of Environmental Quality



**POTENTIAL HAZARDOUS WASTE SITE
SITE INSPECTION REPORT
PART B - OPERATOR INFORMATION**

IDENTIFICATION	
01 STATE	02 SITE NUMBER

WY D980807762

II. CURRENT OPERATOR <small>(Provide if different from owner)</small>		OPERATOR'S PARENT COMPANY <small>(If applicable)</small>			
01 NAME	02 D+B NUMBER	10 NAME	11 D+B NUMBER		
None					
03 STREET ADDRESS <small>(P.O. Box, RFD #, etc.)</small>	04 SIC CODE	12 STREET ADDRESS <small>(P.O. Box, RFD #, etc.)</small>	13 SIC CODE		
05 CITY	06 STATE	07 ZIP CODE	14 CITY	15 STATE	16 ZIP CODE
08 YEARS OF OPERATION	09 NAME OF OWNER				
III. PREVIOUS OPERATOR(S) <small>(List most recent first. Provide only 2 different from owner)</small>		PREVIOUS OPERATORS' PARENT COMPANIES <small>(If applicable)</small>			
01 NAME	02 D+B NUMBER	10 NAME	11 D+B NUMBER		
Ashland Oil Company					
03 STREET ADDRESS <small>(P.O. Box, RFD #, etc.)</small>	04 SIC CODE	12 STREET ADDRESS <small>(P.O. Box, RFD #, etc.)</small>	13 SIC CODE		
Ashland Drive					
05 CITY	06 STATE	07 ZIP CODE	14 CITY	15 STATE	16 ZIP CODE
Ashland	KY				
08 YEARS OF OPERATION	09 NAME OF OWNER DURING THIS PERIOD				
<1	Ashland Oil				
01 NAME	02 D+B NUMBER	10 NAME	11 D+B NUMBER		
Empire State Oil Company					
03 STREET ADDRESS <small>(P.O. Box, RFD #, etc.)</small>	04 SIC CODE	12 STREET ADDRESS <small>(P.O. Box, RFD #, etc.)</small>	13 SIC CODE		
unknown (out of buisness)					
05 CITY	06 STATE	07 ZIP CODE	14 CITY	15 STATE	16 ZIP CODE
08 YEARS OF OPERATION	09 NAME OF OWNER DURING THIS PERIOD				
31	Empire State Oil Company				
01 NAME	02 D+B NUMBER	10 NAME	11 D+B NUMBER		
03 STREET ADDRESS <small>(P.O. Box, RFD #, etc.)</small>	04 SIC CODE	12 STREET ADDRESS <small>(P.O. Box, RFD #, etc.)</small>	13 SIC CODE		
05 CITY	06 STATE	07 ZIP CODE	14 CITY	15 STATE	16 ZIP CODE
08 YEARS OF OPERATION	09 NAME OF OWNER DURING THIS PERIOD				
IV. SOURCES OF INFORMATION <small>(Can include references, e.g., lab sample analysis reports)</small>					
Preliminary Assessment, completed 5/18/84 by Wyoming Department of Environmental Quality Site visit, conducted by Region VIII FIT 6/25/84					

 POTENTIAL HAZARDOUS WASTE SITE SITE INSPECTION REPORT PART 9 - GENERATOR/TRANSPORTER INFORMATION				I. IDENTIFICATION D1 STATE D2 SITE NUMBER WY D980807762	
II. ON-SITE GENERATOR					
D1 NAME None	D2 D+B NUMBER				
D3 STREET ADDRESS (P.O. Box, RFD #, etc.)	D4 SIC CODE				
D5 CITY	D6 STATE	D7 ZIP CODE			
III. OFF-SITE GENERATOR(S)					
D1 NAME None	D2 D+B NUMBER		D1 NAME	D2 D+B NUMBER	
D3 STREET ADDRESS (P.O. Box, RFD #, etc.)	D4 SIC CODE		D3 STREET ADDRESS (P.O. Box, RFD #, etc.)	D4 SIC CODE	
D5 CITY	D6 STATE	D7 ZIP CODE	D5 CITY	D6 STATE	D7 ZIP CODE
D1 NAME None	D2 D+B NUMBER		D1 NAME	D2 D+B NUMBER	
D3 STREET ADDRESS (P.O. Box, RFD #, etc.)	D4 SIC CODE		D3 STREET ADDRESS (P.O. Box, RFD #, etc.)	D4 SIC CODE	
D5 CITY	D6 STATE	D7 ZIP CODE	D5 CITY	D6 STATE	D7 ZIP CODE
IV. TRANSPORTER(S)					
D1 NAME None	D2 D+B NUMBER		D1 NAME	D2 D+B NUMBER	
D3 STREET ADDRESS (P.O. Box, RFD #, etc.)	D4 SIC CODE		D3 STREET ADDRESS (P.O. Box, RFD #, etc.)	D4 SIC CODE	
D5 CITY	D6 STATE	D7 ZIP CODE	D5 CITY	D6 STATE	D7 ZIP CODE
D1 NAME None	D2 D+B NUMBER		D1 NAME	D2 D+B NUMBER	
D3 STREET ADDRESS (P.O. Box, RFD #, etc.)	D4 SIC CODE		D3 STREET ADDRESS (P.O. Box, RFD #, etc.)	D4 SIC CODE	
D5 CITY	D6 STATE	D7 ZIP CODE	D5 CITY	D6 STATE	D7 ZIP CODE
V. SOURCES OF INFORMATION (Give specific references, e.g., sample data, sample analysis, reports.)					
Preliminary Assessment, completed 5/18/84, by Wyoming Department of Environmental Quality Site Visit, conducted by Region VII FIT					



POTENTIAL HAZARDOUS WASTE SITE
SITE INSPECTION REPORT
PART 10 - PAST RESPONSE ACTIVITIES

L IDENTIFICATION

01 STATE	02 SITE NUMBER
WY	D980807762

II. PAST RESPONSE ACTIVITIES

01 A. WATER SUPPLY CLOSED 02 DATE _____ 03 AGENCY _____
04 DESCRIPTION

None

01 B. TEMPORARY WATER SUPPLY PROVIDED 02 DATE _____ 03 AGENCY _____
04 DESCRIPTION

None

01 C. PERMANENT WATER SUPPLY PROVIDED 02 DATE _____ 03 AGENCY _____
04 DESCRIPTION

None

01 D. SPILLED MATERIAL REMOVED 02 DATE _____ 03 AGENCY _____
04 DESCRIPTION

None

01 E. CONTAMINATED SOIL REMOVED 02 DATE _____ 03 AGENCY _____
04 DESCRIPTION

None

01 F. WASTE REPACKAGED 02 DATE _____ 03 AGENCY _____
04 DESCRIPTION

None

01 G. WASTE DISPOSED ELSEWHERE 02 DATE _____ 03 AGENCY _____
04 DESCRIPTION

None

01 H. ON SITE BURIAL 02 DATE _____ 03 AGENCY _____
04 DESCRIPTION

None

01 I. IN SITU CHEMICAL TREATMENT 02 DATE _____ 03 AGENCY _____
04 DESCRIPTION

None

01 J. IN SITU BIOLOGICAL TREATMENT 02 DATE _____ 03 AGENCY _____
04 DESCRIPTION

None

01 K. IN SITU PHYSICAL TREATMENT 02 DATE _____ 03 AGENCY _____
04 DESCRIPTION

None

01 L. ENCAPSULATION 02 DATE _____ 03 AGENCY _____
04 DESCRIPTION

None

01 M. EMERGENCY WASTE TREATMENT 02 DATE _____ 03 AGENCY _____
04 DESCRIPTION

None

01 N. CUTOFF WALLS 02 DATE _____ 03 AGENCY _____
04 DESCRIPTION

None

01 O. EMERGENCY DIKING/SURFACE WATER DIVERSION 02 DATE _____ 03 AGENCY _____
04 DESCRIPTION

None

01 P. CUTOFF TRENCHES/SUMP 02 DATE _____ 03 AGENCY _____
04 DESCRIPTION

None

01 Q. SUBSURFACE CUTOFF WALL 02 DATE _____ 03 AGENCY _____
04 DESCRIPTION

None



POTENTIAL HAZARDOUS WASTE SITE
SITE INSPECTION REPORT
PART 10 - PAST RESPONSE ACTIVITIES

I. IDENTIFICATION	
01 STATE	02 SITE NUMBER
WY	D980807762

II PAST RESPONSE ACTIVITIES (continued)

01 <input type="checkbox"/> R. BARRIER WALLS CONSTRUCTED 04 DESCRIPTION	02 DATE _____	03 AGENCY _____
None		
01 <input type="checkbox"/> S. CAPPING/COVERING 04 DESCRIPTION	02 DATE _____	03 AGENCY _____
None		
01 <input type="checkbox"/> T. BULK TANKAGE REPAIRED 04 DESCRIPTION	02 DATE _____	03 AGENCY _____
None		
01 <input type="checkbox"/> U. GROUT CURTAIN CONSTRUCTED 04 DESCRIPTION	02 DATE _____	03 AGENCY _____
None		
01 <input type="checkbox"/> V. BOTTOM SEALED 04 DESCRIPTION	02 DATE _____	03 AGENCY _____
None		
01 <input type="checkbox"/> W. GAS CONTROL 04 DESCRIPTION	02 DATE _____	03 AGENCY _____
None		
01 <input type="checkbox"/> X. FIRE CONTROL 04 DESCRIPTION	02 DATE _____	03 AGENCY _____
None		
01 <input type="checkbox"/> Y. LEACHATE TREATMENT 04 DESCRIPTION	02 DATE _____	03 AGENCY _____
None		
01 <input type="checkbox"/> Z. AREA EVACUATED 04 DESCRIPTION	02 DATE _____	03 AGENCY _____
None		
01 <input type="checkbox"/> 1. ACCESS TO SITE RESTRICTED 04 DESCRIPTION	02 DATE _____	03 AGENCY _____
None		
01 <input type="checkbox"/> 2. POPULATION RELOCATED 04 DESCRIPTION	02 DATE _____	03 AGENCY _____
None		
01 <input type="checkbox"/> 3. OTHER REMEDIAL ACTIVITIES 04 DESCRIPTION	02 DATE _____	03 AGENCY _____
None		

III. SOURCES OF INFORMATION (Give specific references, e.g., state files, sample analysis, reports.)

Wyoming Department of Environmental Quality files



POTENTIAL HAZARDOUS WASTE SITE
SITE INSPECTION REPORT
PART 11 - ENFORCEMENT INFORMATION

I. IDENTIFICATION	
01 STATE WY	02 SITE NUMBER D980807762

II. ENFORCEMENT INFORMATION

01 PAST REGULATORY/ENFORCEMENT ACTION YES NO

02 DESCRIPTION OF FEDERAL, STATE, LOCAL REGULATORY/ENFORCEMENT ACTION

III. SOURCES OF INFORMATION (List specific references, e.g., state files, sample analysis, reports)

Wyoming Department of Environmental Quality